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U.S. Army Center for Health Promotion and Preventive Medicine

PYROTECHNICS HEALTH RISK ASSESSMENT NO. 39-EJ-1485-99

RESIDENTIAL EXPOSURE FROM INHALATION OF AIR EMISSIONS FROM THE M126A1 RED STAR PARACHUTE SIGNAL FLARE DEPARTMENT OF DEFENSE IDENTIFICATION CODE: L311



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U.S. Army Center for Health Promotion and Preventive Medicine

The lineage of the U.S. Army Center for Health Promotion and Preventive Medicine (USACHPPM) can be traced back over 50 years. This organization began as the U.S. Army Industrial Hygiene Laboratory, established during the industrial buildup for World War II, under the direct supervision of the Army Surgeon General. Its original location was at the Johns Hopkins School of Hygiene and Public Health. Its mission was to conduct occupational health surveys and investigations within the Department of Defense's (DOD's) industrial production base. It was staffed with three personnel and had a limited annual operating budget of three thousand dollars.

Most recently, it became internationally known as the U.S. Army Environmental Hygiene Agency (AEHA). Its mission expanded to support worldwide preventive medicine programs of the Army, DOD, and other Federal agencies as directed by the Army Medical Command or the Office of The Surgeon General, through consultations, support services, investigations, on-site visits, and training.

On 1 August 1994, AEHA was redesignated the U.S. Army Center for Health Promotion and Preventive Medicine with a provisional status and a commanding general officer. On 1 October 1995, the nonprovisional status was approved with a mission of providing preventive medicine and health promotion leadership, direction, and services for America's Army.

The organization's quest has always been one of excellence and the provision of quality service. Today, its goal is to be an established world-class center of excellence for achieving and maintaining a fit, healthy, and ready force. To achieve that end, the CHPPM holds firmly to its values which are steeped in rich military heritage:

- ★ Integrity is the foundation
- ★ Excellence is the standard
- ★ Customer satisfaction is the focus
- ★ Its people are the most valued resource
- ★ Continuous quality improvement is the pathway

This organization stands on the threshold of even greater challenges and responsibilities. It has been reorganized and reengineered to support the Army of the future. The CHPPM now has three direct support activities located in Fort Meade, Maryland; Fort McPherson, Georgia; and Fitzsimons Army Medical Center, Aurora, Colorado; to provide responsive regional health promotion and preventive medicine support across the U.S. There are also two CHPPM overseas commands in Landstuhl, Germany and Camp Zama, Japan who contribute to the success of CHPPM's increasing global mission. As CHPPM moves into the 21st Century, new programs relating to fitness, health promotion, wellness, and disease surveillance are being added. As always, CHPPM stands firm in its commitment to Army readiness. It is an organization proud of its fine history, yet equally excited about its challenging future.

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REPLY TO
ATTENTION OF

MCHB-TS-EHR

PYROTECHNICS HEALTH RISK ASSESSMENT NO. 39-EJ-1485-99
RESIDENTIAL EXPOSURE FROM INHALATION OF AIR EMISSIONS
FROM THE M126A1 RED STAR PARACHUTE SIGNAL FLARE

EXECUTIVE SUMMARY

This assessment evaluated the potential for human health effects to offsite residents breathing air emissions following use of the M126A1 Red star parachute signal flare (M126A1) during training exercises. The military uses pyrotechnics for signaling, obscuring, and illuminating during training and combat. Pyrotechnics are also used during training exercises to simulate battle conditions. Study results showed that no adverse health impacts are expected, to the hypothetical resident, from inhalation of the air emissions from the M126A1.

To conduct this study, air emissions from the M126A1 were collected in a test chamber (Bang Box) at the Dugway Proving Ground, Utah. This information was then used in an air dispersion model to determine ambient air concentrations at a location 100 meters (328 feet) downwind from the site where the M126A1 was activated. Since the training facility in this study is hypothetical, the air model used assumptions that provided conservative estimates of air concentrations.

Modeled air concentrations were combined with exposure information (e.g., number of exposures per year) to estimate the amount of substances the hypothetical resident breathes. This intake was combined with the substance's health information, which was obtained from agencies such as the U.S. Environmental Protection Agency, to determine if there is a potential for health risks from inhalation of these substances.

The health risk study included both long-term (30 years) and short-term (15-minute or 1-hour) exposures to modeled substance concentrations. Study results showed no potential for health risks to the hypothetical resident from inhalation of air emissions from the M126A1.

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LIST OF ACRONYMS

AEC	U.S. Army Environmental Center
AEGL	Acute Exposure Guideline Levels
AIHA	American Industrial Hygiene Association
ATV	Acute Toxicity Value
DODIC	Department of Defense Identification Code
DOD	Department of Defense
DOE	U.S. Department of Energy
EPA	U.S. Environmental Protection Agency
ERPG	Emergency Response Planning Guidelines
HBSL	Health-Based Screening Level
NAAQS	National Ambient Air Quality Standards
NAC/AEGL	National Advisory Committee for Acute Exposure Guideline Levels
NEW	Net Explosive Weight
OEL	Occupational Exposure Limit
PM ₁₀	Particulate Matter under 10 micrometers in size
PRG	Preliminary Remediation Goals
RBC	Risk-Based Concentration
RfC	Reference Concentration
TEEL	Temporary Emergency Exposure Limits
TPCWG	Total Petroleum Criteria Working Group
TSP	Total Suspended Particulates
USACHPPM	U.S. Army Center for Health Promotion and Preventive Medicine

PYROTECHNICS HEALTH RISK ASSESSMENT
NO. 39-EJ-1485-99
RESIDENTIAL EXPOSURE FROM INHALATION OF AIR EMISSIONS FROM THE
M126A1 RED STAR PARACHUTE SIGNAL FLARE

1. PURPOSE

This document presents the evaluation of the potential for human health effects to offsite residents breathing air emissions following use of the M126A1 Red star parachute signal flare (M126A1) during training exercises.

2. AUTHORITY

Memorandum, U.S. Army Environmental Center, 4 June 1999, Subject:
Pyrotechnics Risk Assessment.

3. REFERENCES

See Appendix A.

4. BACKGROUND

a. PYROTECHNICS AND THEIR USE

The term pyrotechnic is derived from the Greek words "pyr" and "techne" meaning fire and art. The terms pyrotechnics and fireworks are often used interchangeably. Examples of pyrotechnics include distress flares and fireworks used for commercial (for public displays) and consumer (e.g., sparklers) use. Every year during New Year and Independence Day celebrations fireworks are used for public displays across the country. During the 1998 Olympic Winter games in Nagano, Japan, almost 5000 pyrotechnics were launched during a firework display that lasted 8 minutes.

The military uses pyrotechnics for four purposes: 1) as a method of communication through the use of signals, 2) to produce smoke to reduce enemy effectiveness, 3) for illuminating the field, and 4) to simulate battle conditions during training exercises. Pyrotechnics play an important role in both military training and combat. It is important that our troops are adequately trained to use them properly.

b. WHAT IS THE M126A1 RED STAR PARACHUTE SIGNAL FLARE?

The M126A1 is a star parachute signal flare, which is a type of pyrotechnic device used for signaling and illuminating. The M126A1 produces a single,

red, parachute-suspended illuminating star. It is 10.16 inches long, 1.67 inches wide, and weighs 1.20 pounds (Reference 1).

c. USE OF THE M126A1 RED STAR PARACHUTE SIGNAL FLARE

The M126A1 is used during many Army training events. These events are held at nearly every Army training installation. In general, one item is used during a day of training, which typically occurs five times per year. A rocket containing the signal is launched from a hand-held device. After igniting, the rocket reaches a height of about 200 feet and produces a single, red star illumination resembling a firework. The signal extends to a height of 700 to 750 feet and can be seen from a distance of 30 to 35 miles at night (References 2, 3). Use of this device is important in training our troops to use and identify different signals, which is an important method of communication in the field.

d. ASSESSMENT SUMMARY

The approach for this study consisted of two main parts: air dispersion modeling and exposure assessment. These are briefly discussed in the paragraphs below. Sections 5-7 present a more explicit discussion of the methodology used for this study.

Data generated in the "Bang Box" at the Dugway Proving Ground, Utah (Reference 4), were used with an atmospheric dispersion model (Reference 5) to estimate the average concentrations that would be experienced by an offsite resident. Since this study is designed to provide results that would be applicable to most Army training facilities, the training area used in this evaluation was a hypothetical one. In addition, air-modeling parameters were selected to mimic worst-case conditions.

The exposure assessment included calculations of time-averaged concentrations for both long-term (chronic) and short-term (acute) exposures. For the purpose of this study, air concentrations were averaged over 30 years for chronic exposures and 1 hour or 15 minutes for acute exposures. These concentrations were compared to chronic health-based screening levels (HBSLs) established by the U.S. Environmental Protection Agency (EPA) or acute toxicity values (ATVs) established by selected agencies depending on the exposure duration (i.e., 30 years versus 1 hour or 15 minutes). If the chronic and acute averaged concentrations (C_{chronic} and C_{acute}) were greater than these screening levels, further analysis would be warranted to determine the potential for health effects. It should be noted that concentrations greater than the screening levels do not indicate an onset of health effects, but rather the potential for such.

5. METHODS AND DATA COLLECTION

a. EMISSION FACTORS

The air modeling emission rates were derived from the pyrotechnics emission studies conducted at Dugway Proving Ground, Utah (Reference 4). These studies sampled air emissions from the firing of weapons and/or munitions used in training. The purpose of this sampling was to identify and quantify the air emissions. The data provided by Dugway Proving Ground included the identification of the munitions item and compounds sampled, net explosive weight (NEW) of the item, vertical and horizontal dimensions of the plume from thermograph data and video, and compound emission factors. This data is included in the tables in Appendix B.

b. AIR MODEL

(1) BACKGROUND

Air dispersion models are available to mathematically simulate atmospheric conditions and behavior to predict downwind concentrations caused by emissions from various sources. However, specific models are not available to estimate the dispersion of emissions from the use of munitions in training. The emissions from munitions used in training result in ambient concentrations of compounds at various locations. The magnitude and location of these concentrations depend on many factors including the amount and type of emissions, the behavior of the source, and meteorological conditions. Based on the evaluation of air dispersion models for military munitions, the U.S. Army Center for Health Promotion and Preventive Medicine (USACHPPM) recommended using the Integrated PUFF (INPUFF) model to estimate the dispersion of emissions from mobile pyrotechnics (Reference 6).

(2) MODEL DESCRIPTION

The INPUFF Model was developed to simulate dispersion from instantaneous or semi-continuous point sources. This Gaussian-integrated puff model is capable of addressing a puff type release over short periods of time, and computations can be performed for a single point source for multiple receptors. The algorithm used to calculate concentrations uses a vertically uniformed wind direction (with no chemical reaction) to compute the contribution of each puff at a receptor for each time step/interval.

(3) ASSUMPTIONS

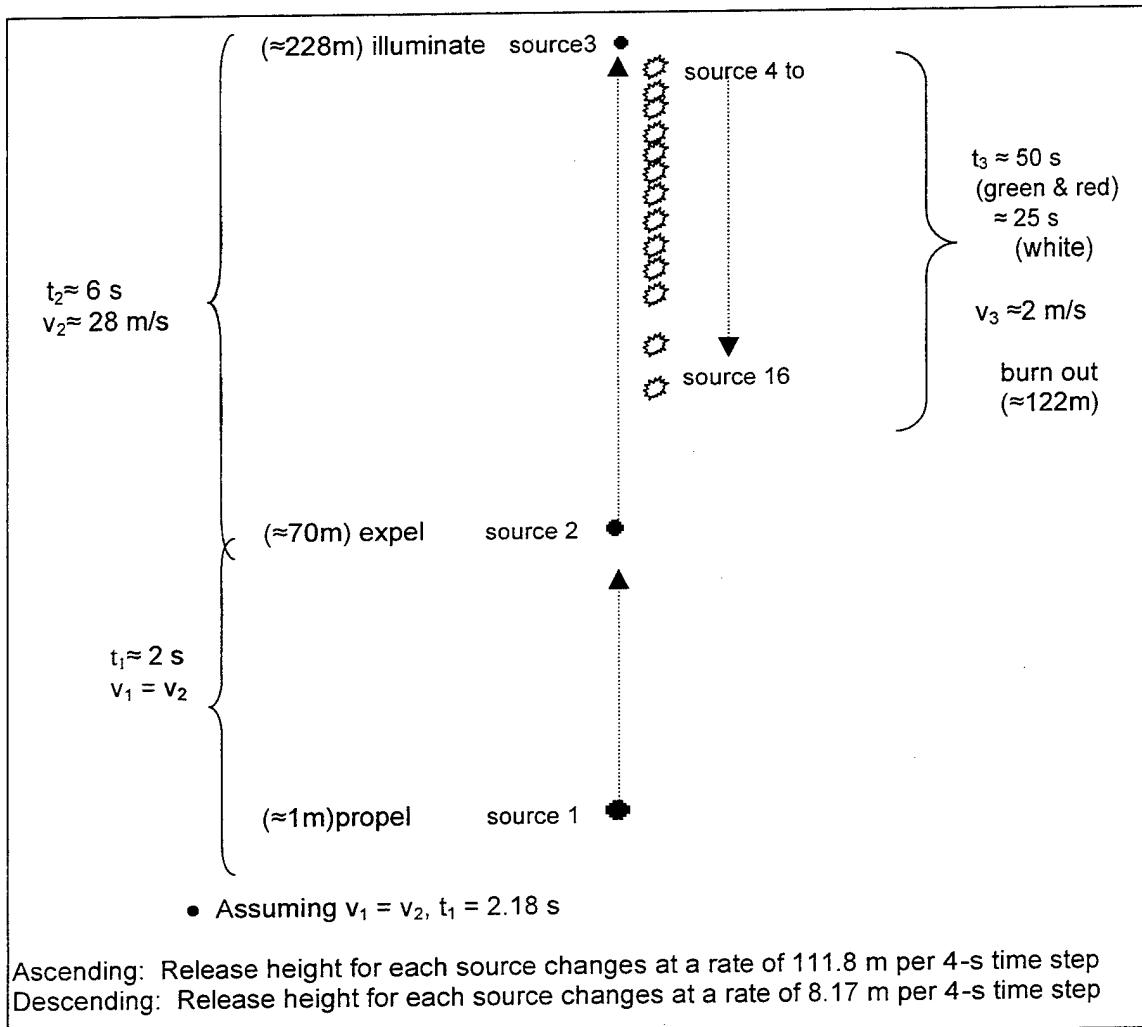
Some assumptions were made to best represent the M126A1 in the air model. These assumptions were as follows:

- (a) For unconventional sources with no physical stack dimensions, the initial horizontal and vertical dispersion values (σ_y and σ_z) of the released puff were used to define the dimensions of the puff. Therefore, plume rise and formation were not determined by characterizing flue gas exit velocity and stack diameter, as they are with conventional point sources. The initial dimensions were set to values measured during Dugway Proving Ground testing and the dispersion of the initial cloud was modeled. The physical dimensions, including height and length of the puff or cloud, were estimated from the thermograph data recorded at every time step. The data also included minimum, mean, and maximum temperature readings during the duration of the emission test and were used to define the flue gas exit temperature. These puffs were then modeled at different release heights as depicted in Figure 1.
- (b) The worst-case release scenario analysis was performed using EPA Risk Management Program Guidance (Reference 7). This guidance includes tables for estimating the footprint of chemical releases. These guidelines were intended to inform emergency responders of the worst possible accidental release, but not necessarily the most likely. The EPA has defined most default conditions for meteorological modeling parameters. Table 1 lists the parameters that were used in the model.
- (c) The resident used in this study was assumed to be directly downwind from the source. The meander of the puff is a major factor when estimating concentrations at given locations downwind from the source. Assuming that the resident is directly downwind from the source is the same as assuming that there is no puff meander and provides the most conservative modeled concentrations.
- (d) For the purpose of this study, the number of items per event was defined as the activation of one item during a 24-hour period.
- (e) Figure 1 provides a schematic diagram of the trajectory of the M126A1. The data for Figure 1 were obtained from References 1 and 8. The performance data provided estimated travel times (t), and velocities (v) at which the items ascend and descend.

TABLE 1. AIR MODEL INPUT PARAMETERS

MODEL PARAMETERS	
Number of meteorological periods (NTIME)	1
Duration of each meteorological period (ITIME)	400 s
Number of updates to the source (NSRCDS)	100 per source
Duration between each source update/time-step (ISUPDT)	4 s
Total time modeled/Simulation Period (NTIME * ITIME)	400 s
SOURCE PARAMETERS	
Source/Stack Diameter	0.5 m for sources (1-16).
Source/Stack Height	See Table 3.
Source Exit Temperature	Varied for each source every time step (4 s) degrees Kelvin (K)
Exit Velocity	NA
Emission Rate	1 g/s
Initial horizontal dispersion (σ_y)	Varied for each source every time step (4 s)
Initial vertical dispersion (σ_z)	Varied for each source every time step (4 s)
WORST CASE METEOROLOGICAL PARAMETERS	
Wind Speed	1 m/s
Atmospheric Stability	Category F
Wind Direction	270° degrees West
Ambient Temperature	293 degrees Kelvin (K) or 68 °F
Worst case Receptor Location	100 m directly downwind

FIGURE 1: SCHEMATIC DIAGRAM OF THE TRAJECTORY OF THE M126A1



(4) GENERAL METHODOLOGY

(a) The INPUFF model determined the amount of time it would take for the puff to pass over a location 100 meters (m) downwind. The released puff migrated at the constant wind speed of one meter per second (1 m/s) downwind from the point of activation. Assuming a distance of 100 m and a travel velocity of 1 m/s, it took 100 s for the center of each puff to reach this distance. The source was modeled at the appropriate release heights and intermediate concentrations were calculated by INPUFF at a receptor located at 100 meters.

(b) The model was run with a total runtime of 400 s to ensure that the total mass of the puff had passed the receptor and the source behavior recorded in the thermograph data was sufficiently simulated. Since the model is capable of providing 100 updates (puffs), each intermediate puff was assumed to be 400 s divided by 100 updates, which is 4 s. Calculated concentrations every time step (4 s) indicated that the puff reached the receptor within 80 s and dissipated below a unit concentration of 1×10^{-10} grams/m³ within 136 s.

(c) The parachute flares were modeled as 16 different sources with different source parameters for each time step. A different release height with varying release temperatures was used for each source. Table 2 illustrates how the emissions for the M126A1 were separated into 16 different point sources and the emissions for each source were staggered over a burn time of 64 s.

TABLE 2: M126A1 SOURCE SIMULATION

Source No.	Height (m)	Unit Emission Rate of 1g/s per time step
1	1.0	0 to 4 s
2	112.0	4 to 8 s
3	228.0	8 to 12 s
4	220.0	12 to 16 s
5	212.0	16 to 20 s
6	203.0	20 to 24 s
7	196.0	24 to 28 s
8	187.0	28 to 32 s
9	179.0	32 to 36 s
10	171.0	36 to 40 s
11	163.0	40 to 44 s
12	154.0	44 to 48 s
13	146.0	48 to 52 s
14	138.0	52 to 56 s
15	130.0	56 to 60 s
16	122.0	60 to 64 s

* The emission rates return to 0 g/s for all time steps not shown in the table. Note: the location of each of the sources with z representing the base elevation was modeled at (x,y,z)=(0,0,0)

(5) USE OF MODEL OUTPUT

The concentrations provided by the INPUFF model are based on a unit emission rate of 1 g/s from an emission source and does not represent any pollutant-specific concentration from the use of pyrotechnics. The relationship between the emission rate and predicted concentration is linear. Therefore, the ratio of the predicted concentration to the unit emission rate was multiplied by each pollutant-specific emission rate to provide pollutant-specific concentrations.

(6) DETERMINATION OF POLLUTANT-SPECIFIC EMISSION RATES

(a) The actual emission rate per item (ER_1) for each pollutant was calculated using the following equation:

$$ER_1 = \frac{M \cdot CV}{t} \quad \text{Equation 1}$$

where:

ER_1 = emission rate for one item (g/(item*sec))
 M = total mass (lb) of pollutant emitted per item (lb/item)
 CV = conversion factor (453.59 g/lb)
 t = release duration in seconds (s) (References 1, 8)

**Example 1
Sample Calculation Using Equation 1*:**

$$ER_1 = \frac{(1.306E - 01)(453.59)}{(64)}$$

$$= 9.254E-01 \text{ g/(s*item)}$$

*Calculation for TSP. Averaged adjusted emission factor of total suspended particulates (TSP) in lb/item was obtained from Appendix B.

(b) The pollutant emission rate for an event (ER_{EV}) for each pollutant was calculated using the estimated number of potential items used in a training event according to the following equation:

$$ER_{EV} = ER_1 \cdot I \quad \text{Equation 2}$$

where:

ER_{EV} = emission rate for the estimated number of potential items used in a training event (g/s)

ER_1 = emission rate for one item (g/(item*sec))

I = items per event (item/event)

Example 2
Sample Calculation Using Equation 2*:

$$ER_{EV} = (9.254E - 01)(1)$$

$$= 9.254E-01 \text{ g/s}$$

* Calculation for TSP

(c) Pollutant-specific ambient concentrations for an event (CONC) were calculated using the following equation:

$$CONC = ER_{EV} \cdot \frac{UC}{ER_{unit}} \quad \text{Equation 3}$$

where:

CONC = pollutant concentration based on the number of items used in a training event (g/m^3)

ER_{EV} = emission rate for the estimated number of items used in a training event (g/s)

ER_{unit} = unit emission rate as used in the model (g/sec)

UC = concentration based on the unit emission rate (g/m^3)

Example 3
Sample Calculation Using Equation 3*:

$$CONC = (9.254E - 01) \frac{(3.028E - 04)}{(1)}$$

$$= 2.802E-04 \text{ g/m}^3$$

* Calculation for TSP

c. EXPOSURE ASSESSMENT

(1) EXPOSURE ASSUMPTIONS

(a) Exposure assumptions were selected using a typical use scenario for the M126A1. This use scenario was developed based on consultation with the U.S. Army Environmental Center's (AEC) senior training advisor (References 9, 10). The frequency of use of the M126A1 was required to determine how much substance an offsite resident would be exposed to in the time period of interest (i.e., acute or chronic exposure). For the purposes of this study, a training scenario is defined as a day or session of training whereas a training event is defined as a single use of pyrotechnics. A training scenario may consist of multiple training events. Table 3 summarizes the specific assumptions used to determine how often the M126A1 is used during a training scenario.

TABLE 3: FREQUENCY OF USE FOR THE M126A1

Parameter	Value Used
Number of items used per training scenario	1
Number of items used per training event	1
Number of training events per scenario	1
Time between events	NA
Number of scenarios per year	5

(b) To estimate the air emissions, it was assumed that one M126A1 was activated. The puff that resulted from this event was modeled to a point 100 meters downwind. Since the unit emission rate was calculated using a runtime of 400 seconds, each event was also assumed to last 400 seconds (or 6.67 minutes).

(2) TIME-AVERAGING

For the chronic assessment, time-averaged concentrations were calculated using the EPA's default residential exposure duration of 30 years (this value assumes that the resident spends 30 years at the same residence). This was done to derive concentrations that would be consistent with the exposure duration used by the EPA so that estimated substance concentrations could be compared to their respective HSLs.

In this evaluation, training scenarios occur five times a year (References 9, 10). Using the default residence time established by the EPA, the assumption was made that someone could be exposed to five training scenarios per year for 30 years.

(a) The daily averaged concentrations were calculated using Equation 4. An example calculation using cadmium is included in Example 4. It should be noted that the average modeled concentration was converted from g/m³ to µg/m³ before it was used in Equation 4.

$$C_d = \frac{CONC \cdot ET \cdot EF_{day}}{1440} \quad \text{Equation 4}$$

where:

C_d = average daily concentration (µg/m³)
CONC = average modeled concentration (µg/m³)
 ET = exposure time (minutes/event)
 EF_{day} = exposure frequency (events/day)
1440 = unit conversion from minutes to day

Example 4
Sample Calculation Using Equation 4:

$$C_{d(cadmium)} = \frac{(1.418E - 03)(6.667)(1)}{1440}$$
$$= 6.565E-06 \mu\text{g}/\text{m}^3$$

The averaged modeled concentration (CONC) for cadmium was obtained from Appendix B. The exposure parameters were obtained from Table 4.

(b) Chronic averaged concentrations were calculated using Equation 5. The resulting concentration (C_d) from Equation 4 was used in Equation 5 to determine the averaged chronic concentrations. Example 5 shows how this calculation was performed.

$$C_{chronic} = \frac{C_d \cdot EF_{year} \cdot ED}{AT} \quad \text{Equation 5}$$

where:

$C_{chronic}$ = average chronic concentration (µg/m³)
 C_d = average daily concentration (µg/m³)
 EF_{year} = exposure frequency (days/year)
 ED = exposure duration (years)
 AT = averaging time (days)
(for carcinogenic endpoint, AT = 70 years x 365 days;
noncarcinogenic endpoint, AT = ED x 365 days)

Example 5
Sample Calculation Using Equation 5:

$$C_{\text{chronic (cadmium)}} = \frac{(6.565E-06)(5)(30)}{(70)(365)}$$

$$= 3.854E-08 \mu\text{g}/\text{m}^3$$

The average daily concentration was calculated as shown in Example 4. The exposure parameters were obtained from Table 4. The averaging time for cadmium is based on the carcinogenic endpoint.

- (c) This study assumed that the same person would be exposed 5 days every year for 30 years. Table 4 lists the exposure parameters used in Equations 4 and 5.

TABLE 4: EXPOSURE PARAMETERS USED TO DETERMINE TIME-AVERAGED CHRONIC AIR CONCENTRATIONS

Exposure Parameter	Value Used
Exposure Time (ET)	6.67 minutes/event
Exposure Frequency (EF _{day})	1 event/day
Exposure Frequency (EF _{year})	5 days/year
Exposure Duration (ED)	30 years

- (d) Unlike the chronic evaluation, guidance for evaluating acute exposures is not currently available. Due to the nature of the use of pyrotechnics and short duration of the concentration plume, acute exposures cannot be overlooked. For the purpose of this study, acute exposure is defined as a 1-hour or 15 minute exposure. The 1-hour or 15 minute acute exposure averaging times allow for comparison with guidelines developed specifically for emergency planning purposes (see discussion on acute toxicity below). This is a conservative assumption since the air model indicated that the hypothetical resident is not expected to be exposed for more than 7 minutes to the concentration plume following activation of the M126A1.
- (e) The average acute concentrations were computed using Equation 6. The exposure frequency is based on the number of events per 1-hour or 15 minutes depending on the guideline used for comparison. Example 6 contains a sample calculation of this equation.

$$C_{acute} = \frac{CONC \cdot ET \cdot EF_{hour}}{60} \quad \text{Equation 6}$$

where:

C_{acute} = average acute concentration ($\mu\text{g}/\text{m}^3$)
CONC = average modeled concentration ($\mu\text{g}/\text{m}^3$)
ET = exposure time (minutes/event)
 EF_{hour} = exposure frequency (events/hour)
60 = unit conversion, 60 minutes/hour

Example 6
Sample Calculation Using Equation 6:

$$C_{acute(cadmium)} = \frac{(1.418E - 03)(6.667)(1 / 0.25)}{60}$$
$$= 6.302E-04 \mu\text{g}/\text{m}^3$$

The average modeled concentration (CONC) for cadmium was obtained from Appendix B. Since the acute toxicity value for cadmium is based on a 15-minute exposure duration (TEEL), the acute concentration was averaged over 15 minutes so that C_{acute} can be compared with its toxicity value.

d. TOXICITY ASSESSMENT

The potential for health risks was determined by comparing time-averaged air concentrations to health-based screening levels, which are developed from a substance's known toxicity. These toxicity values typically include different levels of safety factors depending on the level of confidence of the critical study. Appendix C contains a table of screening values used for the chronic and acute evaluations.

(1) CHRONIC ASSESSMENT

(a) The chronic assessment was evaluated using a screening approach. Using this method, a substance's estimated time-averaged air concentration was compared to its HBSL. If this ratio was less than one, no further analysis was required. This approach is conservative because the exposure assumptions used by the EPA, to establish HBSLs, assume that the resident is exposed for 350 days per year (assuming 2 weeks vacation per year). Since the training scenarios, in which the M126A1 is used, are not expected to exceed 5 days per year, HBSLs specific to this study (if they were developed) would likely be higher.

(b) HBSLs were obtained from the EPA, primarily from Region 3 and Region 9 (References 11, 12). To ensure that the most recent information was used, the Internet sites of both regions were checked. Although the general

approach used by both Region 3 and Region 9 is the same, the exposure assumptions differ enough so that final recommended screening levels can vary to a certain degree. In both methods a substance's HBSL is selected using the toxicity endpoint that derives a lower concentration. For example, if a substance has a known systemic toxicity and is a carcinogen, concentrations were calculated using both toxicity information. The lower concentration was then selected as the recommended screening level to maintain a conservative approach.

(c) A hierarchy was developed in order to quantitatively evaluate for as many of the identified substances as possible. Since the methodology used by Region 9 results in lower HBSLs than Region 3, the Region 9 preliminary remediation goals (PRGs) were used first. Region 3's risk-based concentrations (RBCs) were only used when a PRG was not available. The only exception was for chromium (VI) [Cr (VI)] where Region 9 used a carcinogenic toxicity value that was seven times greater than the EPA's recommended value to develop its screening level for inhalation exposure (Reference 13). Since the EPA does not advocate the application of this multiplication factor, the RBC for Cr (VI) was used instead of the PRG.

(d) Some substances have neither PRGs nor RBCs because they have their own set of regulatory standards. Under the Clean Air Act, the EPA is required to establish National Ambient Air Quality Standards (NAAQS) (Reference 14) for several substances considered harmful to public health and the environment. Currently, NAAQS are available for six substances, of which carbon monoxide, nitrogen dioxide, lead, sulfur dioxide, and particulate matter < 10 micrometers (PM_{10}) have been detected in the M126A1 Bang Box study. The NAAQS for the longer averaging time were used for the chronic evaluation. Depending on the substance, this can range from an 8-hour average to an annual average. In addition, since the majority of the measured total suspended particulates (TSP) were PM_{10} (Reference 4), the NAAQS for PM_{10} was used to evaluate the potential for health effects from exposure to TSP.

Example 7
Sample Calculation Comparing a Substance's Estimated Chronic Concentration to Its Health-Based Screening Level:

$$\frac{C_{chronic(cadmium)}}{HBSL} = \frac{3.85E-08}{1.07E-03}$$
$$= 3.60E-05 < 1$$

The HBSL used for cadmium is a PRG. In this case, the resulting ratio is five orders of magnitude less than one, indicating further evaluation is not necessary.

- (e) Many petroleum hydrocarbons were detected but do not have specific screening levels. Therefore, the approach recommended by the Total Petroleum Criteria Working Group (TPHCWG) (Reference 15) was adopted to evaluate petroleum hydrocarbon mixtures. Based on the working group's assessment of various hydrocarbons, they recommended that mixtures be separated according to a substance's number of carbons and its chemical class (i.e., aliphatic or aromatic¹). Generally, as a substance's carbon number increases, its molecular weight increases and it is therefore, not a substance of concern via inhalation. The working group also concluded that aromatic hydrocarbons tend to be more toxic than aliphatic hydrocarbons (Reference 15).
- (f) Table 5 tabulates the inhalation toxicity values used to evaluate exposure to petroleum mixtures. To be consistent with the methodology used in this study, the reference concentrations (RfCs) were converted to PRGs using Region 9 assumptions. The resulting PRGs are included in Table D-4 in Appendix D.

TABLE 5: SUMMARY OF RfCs USED FOR PETROLEUM HYDROCARBONS (Reference 15)

Carbon Range	Aromatic Inhalation RfC (mg/m ³)	Aliphatic Inhalation RfC (mg/m ³)
C ₅ – C ₆		18.4
C _{>6} – C ₈		
C _{>7} – C ₈	0.4	
C _{>8} – C ₁₀		
C _{>10} – C ₁₂	0.2	1.0
C _{>12} – C ₁₆		
C _{>16} – C ₂₁	NA	NA
C _{>21} – C ₃₅		

NA = not applicable for high molecular weight TPHs (C_{>16}) because compounds in this carbon range are not volatile and therefore, inhalation is not a pathway of concern.

(2) ACUTE ASSESSMENT

- (a) As previously indicated, an acceptable method for assessing acute health effects is not currently available. It was not until recently that EPA guidance addressed the need to evaluate acute health effects from inhalation (Reference 16). Even then, acute toxicity data for risk assessment purposes were not readily available. The EPA recognized this deficiency

¹ Aliphatic hydrocarbons are hydrocarbons in which the carbon atoms are joined by single covalent bonds consisting of two shared electrons (e.g., butane). Aromatic hydrocarbons have ring structures (e.g., benzene) (Reference 21).

and spearheaded the National Advisory Committee for Acute Exposure Guideline Levels for Hazardous Substances (NAC/AEGL Committee). However, AEGLs are currently available for only a handful of substances.

- (b) To circumvent this problem, several state regulatory agencies have suggested that guidelines developed for emergency purposes be used in the interim. Although suggestions have been made to use occupational exposure limits (OELs) by applying additional safety factors (References 17, 18), OELs were not used in this study because they introduce even more uncertainty than the use of emergency guidelines. OELs are designed to protect the workplace environment and assume 8 hours a day, 5 days a week exposures. By definition, these exposures are more chronic than acute.
- (c) In comparison, emergency planning guidelines are more appropriate because they are typically developed for exposures of 1-hour or less. In addition, safety factors may also have been included depending on the agency that develops these guidelines, so that the values would be protective of the general population.
- (d) Emergency Response Planning Guidelines (ERPGs) published by the American Industrial Hygiene Association (AIHA) (Reference 19) and the Temporary Emergency Exposure Limits (TEELs) developed by the U.S. Department of Energy (DOE) (Reference 20) were used for this study; specifically the ERPG-1s and the TEEL-1s. Since TEEL-1s are intended for 15-minute exposures, air concentrations compared to TEELs were averaged over a 15-minute period as opposed to 1-hour in this assessment. The AIHA defines ERPG-1 as follows:

"The maximum concentration in air below which it is believed nearly all individuals could be exposed for up to one hour without experiencing other than mild transient adverse health effects or perceiving a clearly defined objectionable odor."

The DOE defines TEEL-1 as follows:

"The maximum concentration in air below which it is believed nearly all individuals could be exposed without experiencing other than mild transient adverse health effects or perceiving a clearly defined objectionable odor."

(e) For this study, ERPGs were selected prior to a substance's TEEL because they are vigorously reviewed before they are published whereas the TEELs are not. Example 8 shows a sample calculation of how a substance's estimated acute concentration is compared to its acute toxicity value.

Example 8

Sample Calculation of Comparing a Substance's Estimated Acute Concentration to Its Acute Toxicity Value:

$$\frac{C_{\text{acute(cadmium)}}}{ATV} = \frac{6.30E - 04}{3.00E + 01}$$
$$= 2.10E-05 < 1$$

The acute toxicity value available for cadmium is a TEEL. In this example with cadmium, the ratio is five orders of magnitude below 1, indicating that further analysis is not necessary.

6. RISK CHARACTERIZATION

Appendix D presents results from the M126A1 risk characterization. Note that for some substances, two concentrations were reported because of different analytical test methods (as noted in bold). In those instances, the higher concentration was used.

a. CHRONIC HEALTH RISK

The outcome indicated that no chronic health impacts are expected from breathing the air emissions from the M126A1. Since all ratios were below one, no further evaluation was needed.

b. ACUTE HEALTH RISK

For the acute analysis, all ratios were below one, indicating that no acute health impacts are expected from breathing the air emissions from the M126A1. Since all ratios for the acute evaluation were below one, no further assessment was needed.

c. SUBSTANCES WITH NO TOXICITY DATA

Some substances were not quantitatively evaluated because they do not have established toxicity data. By conducting a semi-qualitative comparison of the concentrations of these substances to similar compounds with available toxicity data, it may be concluded that no potential for health effects would be expected from exposure to these substances.

d. FACT SHEET

A copy of the fact sheet submitted to AEC is included as Appendix E. The fact sheet uses the results from this study to summarize health concerns related to inhalation of M126A1 air emissions.

7. UNCERTAINTY DISCUSSION

The limitations inherent in modeling and the added conservatism of the evaluation contribute to the uncertainty of the study results. The risk assessment methodology typically includes safety factors that are embedded in the toxicity data to ensure adequate protection of the general population, particularly, susceptible individuals such as the sick, elderly, and children. Table 6 identifies areas of uncertainty associated with this assessment.

TABLE 6: TYPES OF UNCERTAINTY

Issue	Uncertainty	Direction of Effect
Modeling		
Modeled versus real-time sampling	The air concentrations in this study were modeled. Actual air concentrations taken from the field may be higher or lower.	Varies
Frequency of use for the M126A1	Actual frequency of use of M126A1s during a training event may be different from those stated in this report.	Varies
Hypothetical resident assumed to be located directly downwind	Unless the area around the training facility is populated, the chances that a person living directly downwind is low.	Overestimates
Use of worst-case meteorological conditions	To ensure that this study is applicable to most training areas, worst-case meteorological conditions were used in the air model.	Overestimates
Exposure Assessment		
Estimating time-averaged concentrations	Actual exposure from the M126A1 is intermittent. If one were to plot a person's exposure profile, the plot would consist of a series of spikes. Since current risk assessment methodology does not allow the evaluation of the potential for health risks as a function of time, a single concentration, averaged over the exposure duration was used. In this study, the exposure durations used were 30 years and 1-hour or 15 minutes.	Varies

TABLE 6: TYPES OF UNCERTAINTY

Issue	Uncertainty	Direction of Effect
Chromium speciation	All chromium was assumed to be present as Cr(VI), which is more toxic than Cr(III).	Overestimates
Comparing estimated concentration to established screening levels	The Region 3 and Region 9 HBSLs were developed using different exposure assumptions than those in this study, resulting in more conservative screening levels.	Overestimates
Screening assessment versus calculating an average daily intake	Calculating an average daily intake allows the use of scenario-specific assumptions. However, unless the ratio of concentration to screening level approaches one, a screening assessment is useful as a first-cut evaluation.	Varies
Exposure to other munitions	Other munitions are typically used during the same training event. These items may contain similar or different substances from those detected in the M126A1.	Underestimates
Toxicity Assessment		
Lack of toxicity data	Some substances were not quantitatively evaluated because they have no known toxicity data.	Underestimates
Modifying and uncertainty factors for toxicity data	Modifying factors and uncertainty factors of varying degree are typically applied to toxicological values. These factors are used to conservatively account for extrapolating from animal studies for human health evaluation, and to conservatively account for variation in human populations.	Overestimates

8. CONCLUSION

Results indicated that residents who live as close as 100 meters directly downwind from training areas are safe from breathing air emissions from the M126A1. It is believed that the assumptions contained in this analysis are conservative enough to be protective of all the population including the sick, elderly, and children.

9. RECOMMENDATIONS

Since the results from this study are intended for a hypothetical training facility, they can vary depending on site-specific conditions. However, because of the conservative assumptions used (e.g., worst-case meteorological conditions, receptor located directly downwind, etc.) it is believed that most site-specific analyses would result in even lower concentrations. Therefore, the results from this evaluation should be applicable to most training facilities unless site-specific conditions vary significantly.

10. POINT OF CONTACT

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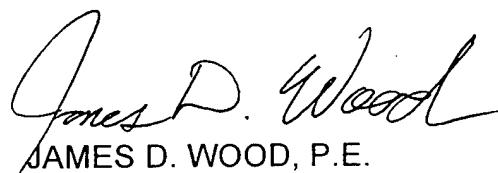
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APPENDIX A

REFERENCES

1. U.S. Army (1994). *Technical Manual, Army Ammunition Data Sheets: Military Pyrotechnics* (Federal Supply Class 1370). TM-43-0001-37.
2. U.S. Army (date unknown). *Chapter 5: Ground Pyrotechnic Signals*. <http://155.217.58.58/cgi-bin/atdl.dll/fm/23-30/Ch5.htm>.
3. U.S. Army (1991). *Operator's Manual, Pyrotechnic Signals*. TM 9-1370-206-10.
4. U.S. Army (1999). *Sampling Results for AEC Phase I Training Ordnance Emission Characterization, Volume I-Summary Report*. Prepared by Radian International LLC, Oak Ridge, TN, for U.S. Army Dugway Proving Ground, Dugway, UT.
5. Bowman Environmental, Inc. (1999). *INPUFF2, Multiple Source Integrated Puff Model*, Version 4.1.
6. USACHPPM (April 2000). *Draft Ambient Air Quality Consultation NO. 43-EL-1485-00 Air Dispersion Modeling Evaluation For Military Munitions*, Aberdeen Proving Ground, MD.
7. Title 40, Code of Federal Regulations, Part 68 (40 CFR 68), Chemical Accident Prevention Provisions, 1 July 1998.
8. U.S. Army Training and Doctrine Command Digital Library, <http://www.adtdl.army.mil/>
9. U.S. Army (1999). Personal communication between Mr. Tony Pitrat, USACHPPM, and Ms. Tamera Clark-Rush, AEC. July 1999.
10. Army Training Evaluation Protocol (ARTEP) 7-20-MTP, *Mission Training Plan for the Infantry Battalion*.
11. EPA (April 1999). *Region 3 Risk Based Concentration (RBC) Tables*. Available online at www.epa.gov/reghwmd/risk/riskmenu.htm
12. EPA (Oct. 1999). *Region 9 Preliminary Remediation Goals (PRG)*. Available online at www.epa.gov/region09/waste/sfund/prg/index.html
13. EPA (1999). *Integrated Risk Information System*. Available online at <http://www.epa.gov/iris/>
14. EPA. *National Ambient Air Quality Standards*. Available online at <http://www.epa.gov/airprog/airs/criteria.html>

15. Total Petroleum Hydrocarbon Criteria Working Group (1997). *Development of Fraction Specific Reference Doses (RfDs) and Reference Concentrations (RfCs) for Total Petroleum Hydrocarbons (TPH)*, Volume 4. Amherst Scientific Publishers. Amherst, MA.
16. EPA (1998). *Human Health Risk Assessment Protocol for Hazardous Waste Combustion Facilities*, Volumes I-III. EPA530-D-98-001A-C.
17. U.S. Army (1996). *Final Screening Risk Assessment for the Anniston Chemical Agent Disposal Facility at the Anniston Army Depot, Alabama*. Revision No. 5. Prepared by the U.S. Army Center for Health Promotion and Preventive Medicine for the Program Manager for Chemical Demilitarization. Aberdeen Proving Ground, Maryland.
18. U.S. Army (1997). *Final Screening Risk Assessment for the Pine Bluff Chemical Agent Disposal Facility at the Pine Bluff Arsenal, Arkansas*. Revision No. 1. Prepared by the U.S. Army Center for Health Promotion and Preventive Medicine for the Program Manager for Chemical Demilitarization. Aberdeen Proving Ground, Maryland.
19. American Industrial Hygiene Association (AIHA). *Emergency Response Planning Guidelines*. AIHA Press, Fairfax, VA.
20. Department of Energy (1998). *Temporary Emergency Exposure Limits*, Revision 15. <http://www.scapa.bnl.gov>.
21. Manahan, Stanley (1994). *Environmental Chemistry*. Sixth edition. CRC Press, Inc. Boca Raton, FL.

APPENDIX B

AIR DISPERSION MODELING OUTPUT DATA

Table B-1: Air Modeling Output Data for Metals, Particulates, and Miscellaneous Compounds

		Red Parachute Signal Flare		Items per event (I):		1 item/event	
		NEW, lb = 0.29		release duration (t):		64 seconds	
		Number of items = 1		Unit Concentration (UC):		3.028E-04 g/m ³ /g(s)	
Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³)	Pollutant Emission Rate (g/sec)/item
				M	CONC	ER ₁	ER _{EV}
Particulate							
TSP	5.739E+01	8.516E-02	4.503E-01	1.306E-01	5.923E+01	2.802E-04	9.254E-01
PM ₁₀	5.735E+01	ND	4.259E-01	1.235E-01	5.602E+01	2.650E-04	8.753E-01
HC Cl ₂							
HCl	2.232E-02	3.676E-02	ND	ND	ND	ND	ND
Cl ₂	6.850E-02	3.646E-02	2.626E-04	7.615E-05	3.454E-02	1.634E-07	5.397E-04
Dioxin Furan							
Dioxin TEQ	3.056E-10	7.123E-11	1.921E-12	5.570E-13	2.527E-10	1.195E-15	3.948E-12
CEM System							
Carbon Monoxide (CO)	4.230E+00	NM (b)	3.594E-02	1.042E-02	4.727E+00	2.237E-05	7.386E-02
Nitrogen Oxide (NOx)	1.272E+00	1.678E-02	1.056E-02	3.063E-03	1.389E+00	6.573E-06	2.171E-02
HCl	3.152E-01	2.725E-01	3.550E-04	1.041E-04	4.722E-02	2.234E-07	7.379E-04
Carbon Dioxide (CO ₂)	7.010E+02	6.429E+02	4.915E-01	1.425E-01	6.465E+01	3.059E-04	1.010E+00
Sulfur Dioxide (SO ₂)	3.279E-02	3.162E-03	2.493E-04	7.231E-05	3.280E-02	1.552E-07	5.125E-04
Particulate-phase Metals							
Aluminum	7.480E-02	NM (a)	5.877E-04	1.704E-04	7.731E-02	3.658E-07	1.208E-03
Antimony	1.958E-04	NM (a)	1.559E-06	4.462E-07	2.024E-04	9.575E-10	3.162E-06
Arsenic	ND	NM (a)	ND	ND	ND	ND	ND
Barium	1.585E-01	NM (a)	1.245E-03	3.611E-04	1.638E-01	7.749E-07	2.559E-03
Beryllium	ND	NM (a)	ND	ND	ND	ND	ND
Cadmium	2.899E-04	NM (a)	2.278E-06	6.606E-07	2.997E-04	1.418E-09	4.682E-06
Chromium	1.505E-03	NM (a)	1.182E-05	3.429E-06	1.555E-03	7.359E-09	2.430E-05
Cobalt	1.967E-04	NM (a)	1.545E-06	4.481E-07	2.033E-04	9.616E-10	3.176E-06
Copper	3.031E-03	NM (a)	2.382E-05	6.906E-06	3.133E-03	1.482E-08	4.895E-05
Lead	1.108E-03	NM (a)	8.704E-06	2.524E-06	1.145E-03	5.417E-09	1.789E-05
Magnesium	6.836E+00	NM (a)	5.372E-02	1.558E-02	7.066E+00	3.343E-05	1.104E-01
Manganese	1.042E-02	NM (a)	8.187E-05	2.374E-05	1.077E-02	5.095E-08	1.683E-04

Table B-1: Air Modeling Output Data for Metals, Particulates, and Miscellaneous Compounds

Compound	Red Parachute Signal Flare		Items per event (I): release duration (t): Unit Concentration (UC):	1 item/event 64 seconds 3.028E-04 g/m ³ /g/s)	
	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)		Pollutant Concentration 1 Item (grams/m ³)	Pollutant Emission Rate (g/sec)/item
		Average Adjusted Emission Factor (lb/lb NEW)			
Nickel	3.014E-04	NM (a)	2.368E-06	6.866E-07	3.115E-04
Phosphorus	5.149E-03	NM (a)	4.046E-05	1.173E-05	5.322E-03
Selenium	ND	NM (a)	ND	ND	2.518E-08
Silver	ND	NM (a)	ND	ND	ND
Thallium	ND	NM (a)	ND	ND	ND
Zinc	2.879E-03	NM (a)	2.263E-05	6.560E-06	2.975E-03
Mercury	4.105E-05	NM (a)	3.226E-07	9.354E-08	4.243E-05

Footnotes:

ND = Not Detected

NEW = Net Explosive Weight

NM = Not Measurable

a = Insufficient material to analyze.

b = Concentration reported was less than zero.

Table B-2: Air Modeling Output Data for Volatile Organic Compounds

Compound	Red Parachute Signal Flare				Items per event (I):		1 item/event	
	NEW, lb = 0.29	Number of Items = 1	release duration (t):	Unit Concentration (UC):	3.028E-04	64 seconds	g/m ³ /sec)	
Total Nonmethane Hydrocarbons (TNMHC)	2.051E-01	5.460E-02	1.153E-03	3.343E-04	1.517E-01	7.175E-07	2.370E-03	2.370E-03
Volatile Organic Compounds (VOCs)								
Ethane	7.150E-03	1.600E-03	4.253E-05	1.233E-05	5.595E-03	2.647E-08	8.741E-05	
Ethylene	5.835E-02	2.000E-04	4.456E-04	1.292E-04	5.862E-02	2.773E-07	9.159E-04	
Acetylene	4.110E-02	6.000E-04	3.104E-04	9.000E-05	4.083E-02	1.932E-07	6.379E-04	
Propane	1.950E-03	7.000E-04	9.579E-06	2.778E-06	1.260E-03	5.962E-09	1.969E-05	
Propene	1.305E-02	3.000E-04	9.771E-05	2.833E-05	1.285E-02	6.081E-08	2.008E-04	
i-Butane	7.500E-04	5.000E-04	1.916E-06	5.556E-07	2.520E-04	1.192E-09	3.938E-06	
j-Butene	7.000E-04	ND	5.364E-06	1.556E-06	7.056E-04	3.338E-09	1.103E-05	
1-Butene	2.500E-03	ND	1.916E-05	5.556E-06	2.520E-03	1.192E-08	3.938E-05	
1,3-Butadiene	3.200E-03	ND	2.452E-05	7.111E-06	3.226E-03	1.526E-08	5.040E-05	
n-Butane	2.150E-03	1.800E-03	2.682E-06	7.778E-07	3.528E-04	1.669E-09	5.513E-06	
trans-2-Butene	2.500E-03	ND	1.916E-05	5.556E-06	2.520E-03	1.192E-08	3.938E-05	
2,2-Dimethylpropane	ND	ND	ND	ND	ND	ND	ND	ND
cis-2-Butene	4.500E-04	ND	3.448E-06	1.000E-06	4.536E-04	2.146E-09	7.088E-06	
3-Methyl-1-butene	ND	ND	ND	ND	ND	ND	ND	ND
i-Pentane	3.100E-03	3.100E-03	ND	ND	ND	ND	ND	ND
1-Pentene	5.000E-04	ND	3.833E-06	1.111E-06	5.040E-04	2.385E-09	7.875E-06	
2-Methyl-1-butene	1.000E-04	ND	7.663E-07	2.222E-07	1.0008E-04	4.769E-10	1.575E-06	
n-Pentane	2.700E-03	2.100E-03	4.598E-06	1.333E-06	6.048E-04	2.862E-09	9.450E-06	
Isoprene	5.000E-04	ND	3.832E-06	1.111E-06	5.040E-04	2.385E-09	7.875E-06	
trans-2-Pentene	1.000E-04	ND	7.663E-07	2.222E-07	1.0008E-04	4.769E-10	1.575E-06	
cis-2-Pentene	1.000E-04	ND	7.663E-07	2.222E-07	1.0008E-04	4.769E-10	1.575E-06	
2-Methyl-2-butene	2.000E-04	ND	1.533E-06	4.445E-07	2.016E-04	9.538E-10	3.150E-06	
2,2-Dimethylbutane	6.000E-04	4.000E-04	1.533E-06	4.445E-07	2.016E-04	9.538E-10	3.150E-06	
Cyclopentene	1.500E-04	ND	1.149E-06	3.333E-07	1.512E-04	7.154E-10	2.363E-06	
4-Methyl-1-pentene	ND	ND	ND	ND	ND	ND	ND	ND
Cyclopentane	1.500E-04	1.000E-04	3.832E-07	1.111E-07	5.040E-05	2.385E-10	7.875E-07	
2,3-Dimethylbutane	2.500E-04	2.000E-04	3.832E-07	1.111E-07	5.040E-05	2.385E-10	7.875E-07	
cis-4-Methyl-2-pentene	ND	ND	ND	ND	ND	ND	ND	ND
2-Methylpentane	1.200E-03	1.100E-03	7.663E-07	2.222E-07	1.0008E-04	4.769E-10	1.575E-06	

Table B-2: Air Modeling Output Data for Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/ item NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³) CONC	Pollutant Emission Rate (g/sec)/item	Event Pollutant Emission Rate 1 Item (g/sec) ER _{EV}
3-Methylpentane	5.000E-04	6.000E-04	ND	ND	ND	ND	ND	ND
2-Methyl-1-pentene	ND	ND	ND	ND	ND	ND	ND	ND
1-Hexene	4.000E-04	ND	3.065E-06	8.889E-07	4.032E-04	1.908E-09	6.300E-06	6.300E-06
n-Hexane	1.100E-03	1.000E-03	7.663E-07	2.222E-07	1.008E-04	4.769E-10	1.575E-06	1.575E-06
trans-2-Hexene	ND	ND	ND	ND	ND	ND	ND	ND
2-Methyl-2-pentene	ND	ND	ND	ND	ND	ND	ND	ND
cis-2-Hexene	ND	ND	ND	ND	ND	ND	ND	ND
Methylcyclopentane	5.500E-04	4.000E-04	1.149E-06	3.333E-07	1.512E-04	7.154E-10	2.363E-06	2.363E-06
2,4-Dimethylpentane	1.000E-04	1.000E-04	ND	ND	ND	ND	ND	ND
Benzene	6.900E-03	1.100E-03	4.445E-05	1.289E-05	5.847E-03	2.766E-08	9.135E-05	9.135E-05
Cyclohexane	4.000E-04	ND	3.065E-06	8.889E-07	4.032E-04	1.908E-09	6.300E-06	6.300E-06
2-Methylhexane	4.000E-04	4.000E-04	ND	ND	ND	ND	ND	ND
2,3-Dimethylpentane	4.000E-04	5.000E-04	ND	ND	ND	ND	ND	ND
3-Methylhexane	4.500E-04	6.000E-04	ND	ND	ND	ND	ND	ND
2,2,4-Trimethylpentane	1.250E-03	1.300E-03	ND	ND	ND	ND	ND	ND
n-Heptane	6.500E-04	4.000E-04	1.916E-06	5.566E-07	2.520E-04	1.192E-09	3.938E-06	3.938E-06
2,4,4-Trimethyl-1-pentene	ND	ND	ND	ND	ND	ND	ND	ND
Methylcyclohexane	3.500E-04	4.000E-04	ND	ND	ND	ND	ND	ND
2,4,4-Trimethyl-2-pentene	ND	ND	ND	ND	ND	ND	ND	ND
2,5-Dimethylhexane	1.500E-04	1.000E-04	3.383E-07	1.111E-07	5.040E-05	2.385E-10	7.875E-07	7.875E-07
2,4-Dimethylhexane	1.000E-04	2.000E-04	ND	ND	ND	ND	ND	ND
2,3,4-Trimethylpentane	3.000E-04	3.000E-04	4.154E-22	1.205E-22	5.465E-20	2.585E-26	8.538E-22	8.538E-22
Toluene	3.700E-03	2.500E-03	9.196E-06	2.667E-06	1.210E-03	5.723E-09	1.890E-05	1.890E-05
2,3-Dimethylhexane	1.000E-04	2.000E-04	ND	ND	ND	ND	ND	ND
2-Methylheptane	2.000E-04	1.000E-04	7.663E-07	2.222E-07	1.008E-04	4.769E-10	1.575E-06	1.575E-06
3-Ethylhexane	2.000E-04	1.000E-04	7.663E-07	2.222E-07	1.008E-04	4.769E-10	1.575E-06	1.575E-06
2,2-Dimethylheptane	ND	ND	ND	ND	ND	ND	ND	ND
2,2,4-Trimethylhexane	ND	1.000E-04	ND	ND	ND	ND	ND	ND
n-Octane	2.000E-04	1.000E-04	7.663E-07	2.222E-07	1.008E-04	4.769E-10	1.575E-06	1.575E-06
Ethylcyclohexane	ND	ND	ND	ND	ND	ND	ND	ND
Ethylbenzene	1.100E-03	1.300E-03	ND	ND	ND	ND	ND	ND
m-Xylene & p-Xylene	4.350E-03	5.400E-03	ND	ND	ND	ND	ND	ND
Styrene	4.500E-04	ND	3.448E-06	1.000E-06	4.536E-04	2.146E-09	7.088E-06	7.088E-06
c-Xylene	1.500E-03	2.000E-03	ND	ND	ND	ND	ND	ND
n-Nonane	7.500E-04	ND	5.747E-06	1.667E-06	7.560E-04	3.577E-09	1.181E-05	1.181E-05
i-Propylbenzene	ND	ND	ND	ND	ND	ND	ND	ND

Table B-2: Air Modeling Output Data for Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³) CONC	Pollutant Emission Rate (g/sec)/item	Pollutant Emission Rate 1 Item (g/sec)	Event Pollutant Emission Rate 1 Item (g/sec) EREV
n-Propylbenzene	2.000E-04	ND	1.533E-06	4.445E-07	2.016E-04	9.538E-10	3.150E-06	3.150E-06	
p-Ethyltoluene	5.000E-04	ND	3.832E-06	1.111E-06	5.040E-04	2.385E-09	7.875E-06	7.875E-06	
m-Ethyltoluene	3.000E-04	ND	2.299E-06	6.667E-07	3.024E-04	1.431E-09	4.725E-06	4.725E-06	
1,3,5-Trimethylbenzene	2.500E-04	ND	1.916E-06	5.556E-07	2.520E-04	1.192E-09	3.938E-06	3.938E-06	
o-Ethyltoluene	4.000E-04	ND	3.065E-06	8.889E-07	4.032E-04	1.908E-09	6.300E-06	6.300E-06	
1,2,4-Trimethylbenzene & sec-Butylbenzene	6.000E-04	ND	4.598E-06	1.333E-06	6.048E-04	2.862E-09	9.450E-06	9.450E-06	
n-Decane	ND	2.000E-04	ND	ND	ND	ND	ND	ND	ND
alpha-Pinene	ND	ND	ND	ND	ND	ND	ND	ND	ND
beta-Pinene	ND	ND	ND	ND	ND	ND	ND	ND	ND
delta-3-Carene	ND	ND	ND	ND	ND	ND	ND	ND	ND
d-Limonene	ND	ND	ND	ND	ND	ND	ND	ND	ND
MTBE	ND	4.000E-04	ND	ND	ND	ND	ND	ND	ND
Dichlorodifluoromethane	1.653E-03	1.680E-03	ND	ND	ND	ND	ND	ND	ND
Methylchloride	ND	ND	ND	ND	ND	ND	ND	ND	ND
Dichlorotetrafluoroethane	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chloroethene	ND	ND	ND	ND	ND	ND	ND	ND	ND
1,3-Butadiene	3.255E-03	ND	2.494E-05	7.233E-06	3.281E-03	1.552E-08	5.126E-05	5.126E-05	
Methylbromide	ND	ND	ND	ND	ND	ND	ND	ND	ND
Ethylchloride	1.398E-04	ND	1.072E-06	3.108E-07	1.410E-04	6.669E-10	2.203E-06	2.203E-06	
Trichloromonofluoromethane	2.650E-03	2.785E-03	ND	ND	ND	ND	ND	ND	
Vinyldenechloride	ND	ND	ND	ND	ND	ND	ND	ND	
Methylenechloride	ND	ND	ND	ND	ND	ND	ND	ND	
Allylchloride	ND	ND	ND	ND	ND	ND	ND	ND	
1,1,2-Trichloro-1,2,2-trifluoroethane	9.017E-04	9.470E-04	ND	ND	ND	ND	ND	ND	
1,1-Dichloroethane	ND	ND	ND	ND	ND	ND	ND	ND	
1,2-Dichloroethene	ND	ND	ND	ND	ND	ND	ND	ND	
Chloroform	ND	ND	ND	ND	ND	ND	ND	ND	
1,2-Dichloroethane	ND	ND	ND	ND	ND	ND	ND	ND	
Methylchloroform	3.200E-04	3.474E-04	ND	ND	ND	ND	ND	ND	
Benzene	7.018E-03	1.119E-03	4.521E-05	1.311E-05	5.947E-03	2.814E-08	9.292E-05	9.292E-05	
Carbon tetrachloride	7.155E-04	8.244E-04	ND	ND	ND	ND	ND	ND	
1,2-Dichloropropane	ND	ND	ND	ND	ND	ND	ND	ND	
Trichloroethylene	ND	ND	ND	ND	ND	ND	ND	ND	
cis 1,3-Dichloro-1-propene	ND	ND	ND	ND	ND	ND	ND	ND	
trans 1,3-Dichloro-1-propene	ND	ND	ND	ND	ND	ND	ND	ND	
1,1,2-Trichloroethane	ND	ND	ND	ND	ND	ND	ND	ND	

Table B-2: Air Modeling Output Data for Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³) CONC	Pollutant Emission Rate 1 Item (g/sec) ER ₁	Event Pollutant Emission Rate 1 Item (g/sec) ER _{EV}
Toluene	3.763E-03	2.543E-03	9.353E-06	2.712E-06	1.230E-03	5.821E-09	1.922E-05	1.922E-05
1,2-Dibromoethane	ND	ND	ND	ND	ND	ND	ND	ND
Perchloroethylene	ND	ND	ND	ND	ND	ND	ND	ND
Chlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
Ethylbenzene	1.689E-03	1.996E-03	ND	ND	ND	ND	ND	ND
m&p-Xylene	4.424E-03	5.492E-03	ND	ND	ND	ND	ND	ND
Styrene	4.577E-04	ND	3.507E-06	1.017E-06	4.614E-04	2.183E-09	7.209E-06	7.209E-06
1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	ND	ND	ND	ND
o-Xylene	1.526E-03	2.034E-03	ND	ND	ND	ND	ND	ND
p-Ethyltoluene	5.086E-04	ND	3.897E-06	1.130E-06	5.126E-04	2.425E-09	8.010E-06	8.010E-06
1,3,5-Trimethylbenzene	3.051E-04	ND	2.338E-06	6.781E-07	3.076E-04	1.455E-09	4.806E-06	4.806E-06
1,2,4-Trimethylbenzene	6.103E-04	ND	4.677E-06	1.356E-06	6.152E-04	2.911E-09	9.612E-06	9.612E-06
Benzylchloride	ND	ND	ND	ND	ND	ND	ND	ND
m-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
p-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
o-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-Trichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
Hexachlorobutadiene	ND	ND	ND	ND	ND	ND	ND	ND
trans-1,2-Dichloroethene	ND	ND	ND	ND	ND	ND	ND	ND
o-Chlorotoluene	ND	ND	ND	ND	ND	ND	ND	ND
p-Chlorotoluene	ND	ND	ND	ND	ND	ND	ND	ND
1,3,5-Trichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
1,2,3-Trichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
Methylnitrite	4.068E-04	3.025E-04	7.992E-07	2.318E-07	1.051E-04	4.974E-10	1.643E-06	1.643E-06
Acetonitrile	4.976E-04	ND	3.873E-06	1.106E-06	5.016E-04	2.373E-09	7.838E-06	7.838E-06
Acrylonitrile	3.813E-04	ND	2.922E-06	8.473E-07	3.843E-04	1.818E-09	6.005E-06	6.005E-06
Nitromethane	7.151E-04	4.677E-04	1.896E-06	5.500E-07	2.495E-04	1.180E-09	3.898E-06	3.898E-06
Benzonitrile	2.149E-04	ND	1.647E-06	4.775E-07	2.166E-04	1.025E-09	3.384E-06	3.384E-06
Nitrobenzene	ND	ND	ND	ND	ND	ND	ND	ND
Carbonyl Sulfide	2.057E-04	ND	1.576E-06	4.570E-07	2.073E-04	9.808E-10	3.239E-06	3.239E-06
Sulfur Dioxide	ND	ND	ND	ND	ND	ND	ND	ND
Carbon Disulfide	6.686E-03	5.214E-04	4.724E-05	1.370E-05	6.214E-03	2.940E-08	9.709E-05	9.709E-05
Thiophene	2.582E-04	ND	1.979E-06	5.739E-07	2.603E-04	1.232E-09	4.067E-06	4.067E-06
Dimethyldisulfide	ND	ND	ND	ND	ND	ND	ND	ND
2-Methylthiophene	ND	ND	ND	ND	ND	ND	ND	ND
3-Methylthiophene	ND	ND	ND	ND	ND	ND	ND	ND

Table B-2: Air Modeling Output Data for Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Concentration 1 Item (grams/m ³) CONC	Pollutant Emission Rate (g/sec)/item ER ₁	Event Pollutant Emission Rate 1 Item (g/sec) ER _{EV}
Dimethyltrisulfide	ND	ND	ND	ND	ND	ND	ND	ND
Isothiocyanatomethane	ND	ND	ND	ND	ND	ND	ND	ND
2-Chlorothiophene	ND	ND	ND	ND	ND	ND	ND	ND
3-Chlorothiophene	ND	ND	ND	ND	ND	ND	ND	ND
2-Thiophenecarboxaldehyde	ND	ND	ND	ND	ND	ND	ND	ND
Naphthalene	3.274E-04	ND	2.509E-06	7.275E-07	3.300E-04	1.561E-09	5.156E-06	5.156E-06
Acetaldehyde	6.358E-04	ND	4.872E-06	1.413E-06	6.409E-04	3.032E-09	1.001E-05	1.001E-05
Acrolein	2.649E-04	ND	2.030E-06	5.888E-07	2.671E-04	1.264E-09	4.173E-06	4.173E-06
Acetone	7.386E-03	8.413E-03	ND	ND	ND	ND	ND	ND
Propanal	6.012E-04	ND	4.607E-06	1.336E-06	6.060E-04	2.867E-09	9.468E-06	9.468E-06
Furan	2.643E-04	ND	2.025E-06	5.874E-07	2.664E-04	1.261E-09	4.163E-06	4.163E-06
2-Propanol	ND	ND	ND	ND	ND	ND	ND	ND
2-Methylpropanol	ND	ND	ND	ND	ND	ND	ND	ND
Methacrolein	ND	ND	ND	ND	ND	ND	ND	ND
2,3-Butanedione	ND	ND	ND	ND	ND	ND	ND	ND
Methyl-Vinyl Ketone	ND	ND	ND	ND	ND	ND	ND	ND
MTBE	8.816E-04	8.413E-04	3.059E-07	8.959E-08	4.064E-05	1.923E-10	6.349E-07	6.349E-07
Butanal	9.798E-04	6.832E-04	2.273E-06	6.591E-07	2.989E-04	1.414E-09	4.671E-06	4.671E-06
2-Butanone	1.253E-03	7.843E-04	3.590E-06	1.041E-06	4.722E-04	2.234E-09	7.378E-06	7.378E-06
Tetrahydrofuran	ND	ND	ND	ND	ND	ND	ND	ND
2-Methyl-1-propanol	ND	ND	ND	ND	ND	ND	ND	ND
trans-2-Butenal	ND	ND	ND	ND	ND	ND	ND	ND
Acetic Acid	1.378E-03	8.749E-04	3.856E-06	1.118E-06	5.073E-04	2.400E-09	7.926E-06	7.926E-06
2-Pentanone	1.642E-03	1.061E-03	4.453E-06	1.291E-06	5.858E-04	2.772E-09	9.153E-06	9.153E-06
Pentanal	1.606E-03	1.560E-03	3.598E-07	1.044E-07	4.733E-05	2.239E-10	7.396E-07	7.396E-07
4-Methyl-2-pentanone	ND	ND	ND	ND	ND	ND	ND	ND
trans-2-Pentenal	ND	ND	ND	ND	ND	ND	ND	ND
Cyclopentanone	ND	ND	ND	ND	ND	ND	ND	ND
2-Hexanone	2.181E-04	ND	1.671E-06	4.847E-07	2.198E-04	1.040E-09	3.435E-06	3.435E-06
Hexanal	5.276E-04	9.664E-04	ND	ND	ND	ND	ND	ND
3-Furaldehyde	ND	ND	ND	ND	ND	ND	ND	ND
Butyl Acetate	ND	ND	ND	ND	ND	ND	ND	ND
2-Furaldehyde	ND	ND	ND	ND	ND	ND	ND	ND
trans-2-Hexenal	ND	ND	ND	ND	ND	ND	ND	ND
1-Hexanol	ND	ND	ND	ND	ND	ND	ND	ND
3-Heptanone	5.014E-04	ND	3.842E-06	1.114E-06	5.054E-04	2.391E-09	7.897E-06	7.897E-06

Table B-2: Air Modeling Output Data for Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³) CONC	Pollutant Emission Rate (g/sec)/item	Event Pollutant Emission Rate 1 item (g/sec)
				M		ER ₁	ER _{REV}	
2-Heptanone	ND	ND	ND	ND	ND	ND	ND	ND
Heptanal	6.160E-04	1.281E-03	ND	ND	ND	ND	ND	ND
trans-2-Heptenal	ND	ND	ND	ND	ND	ND	ND	ND
5-Methyl-2-furaldehyde	ND	ND	ND	ND	ND	ND	ND	ND
6-Methyl-2-heptanone	ND	ND	ND	ND	ND	ND	ND	ND
Benzaldehyde	1.272E-03	1.037E-03	1.798E-06	5.214E-07	2.365E-04	1.119E-09	3.695E-06	3.695E-06
1-Heptanol	ND	ND	ND	ND	ND	ND	ND	ND
6-Methyl-5-hepten-2-one	ND	5.975E-04	ND	ND	ND	ND	ND	ND
2-Octanone	ND	ND	ND	ND	ND	ND	ND	ND
Octanal	1.874E-03	2.740E-03	ND	ND	ND	ND	ND	ND
Benzofuran	ND	ND	ND	ND	ND	ND	ND	ND
trans-2-Octenal	ND	ND	ND	ND	ND	ND	ND	ND
Acetophenone	4.455E-04	3.578E-04	6.720E-07	1.949E-07	8.840E-05	4.182E-10	1.381E-06	1.381E-06
2-Nonanone	ND	ND	ND	ND	ND	ND	ND	ND
Nonanal	3.308E-03	1.519E-03	1.371E-05	3.976E-06	1.803E-03	8.533E-09	2.818E-05	2.818E-05
trans-2-Nonenal	ND	ND	ND	ND	ND	ND	ND	ND
2-Decanone	ND	ND	ND	ND	ND	ND	ND	ND
Decanal	2.662E-03	ND	2.040E-05	5.91626E-06	2.684E-03	1.270E-08	4.193E-05	4.193E-05

Footnotes:

ND = Not Detected

NEW = Net Explosive Weight

Items in bold represent duplicate values for those compounds that are common for Method TO-14 and TO-12.

Table B-3: Air Modeling Output Data for Semi-Volatile Organic Compounds

Compound	Red Parachute Signal Flare		Items per event (I):		1 item/event	
	NEW, lb = 0.29	release duration (t):	64 seconds	Unit Concentration (UC):	3.028E-04 g/m ³ /s	Event Pollutant Emission Rate 1 Item (g/sec)
	Number of Items = 1		Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³)	Pollutant Emission Rate (g/sec)/item	ER _{EV}
	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/lb Item)	M	CONC
Particulate/apor-phase SVOCs						
N-Nitrosodimethylamine	ND	ND	ND	ND	ND	ND
Pyridine	ND	ND	ND	ND	ND	ND
2-Picoline	ND	ND	ND	ND	ND	ND
Methyl methanesulfonate	ND	ND	ND	ND	ND	ND
N-Nitrosomethyl ethylamine	ND	ND	ND	ND	ND	ND
N-Nitrosodiethylamine	ND	ND	ND	ND	ND	ND
Ethyl methanesulfonate	ND	ND	ND	ND	ND	ND
Phenol	ND	ND	ND	ND	ND	ND
Aniline	ND	ND	ND	ND	ND	ND
bis(2-Chloroethyl)ether	ND	ND	ND	ND	ND	ND
Pentachloroethane	ND	ND	ND	ND	ND	ND
2-Chlorophenol	ND	ND	ND	ND	ND	ND
1,3-Dichlorobenzene	ND	ND	ND	ND	ND	ND
1,4-Dichlorobenzene	ND	ND	ND	ND	ND	ND
Benzyl alcohol	ND	ND	ND	ND	ND	ND
2-Methylphenol	ND	ND	ND	ND	ND	ND
1,2-Dichlorobenzene	ND	ND	ND	ND	ND	ND
bis(2-Chloroisopropyl)ether	ND	ND	ND	ND	ND	ND
c-Toluidine	ND	ND	ND	ND	ND	ND
4-Methylphenol/3-Methylphenol	ND	ND	ND	ND	ND	ND
N-Nitroso-di-n-propylamine	ND	ND	ND	ND	ND	ND
Acetophenone	1.279E-03	3.066E-04	7.967E-06	2.310E-06	1.048E-03	4.958E-09
N-Nitrosomorpholine	ND	ND	ND	ND	ND	1.637E-05
N-Nitrosopyrrolidine	ND	ND	ND	ND	ND	ND
Hexachloroethane	ND	ND	ND	ND	ND	ND
Nitrobenzene	ND	ND	ND	ND	ND	ND
N-Nitrosopiperidine	ND	ND	ND	ND	ND	ND
Isophorone	ND	ND	ND	ND	ND	ND
2,4-Dimethylphenol	ND	ND	ND	ND	ND	ND
2-Nitrophenol	ND	ND	ND	ND	ND	ND
bis(2-Chloroethoxy)methane	ND	ND	ND	ND	ND	ND
Benzoic acid	3.493E-03	2.394E-03	9.012E-06	2.613E-06	1.185E-03	5.609E-09
						1.852E-05

Table B-3: Air Modeling Output Data for Semi-Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³)	Pollutant Emission Rate (g/sec)/item	Event Pollutant Emission Rate 1 Item (g/sec)
				M		CONC	ER ₁	ER _{EV}
2,4-Dichlorophenol	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4-Trichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
Naphthalene	ND	ND	ND	ND	ND	ND	ND	ND
p-Chloroaniline	ND	ND	ND	ND	ND	ND	ND	ND
2,6-Dichlorophenol	ND	ND	ND	ND	ND	ND	ND	ND
Hexachloropropene	ND	ND	ND	ND	ND	ND	ND	ND
Hexachlorobutadiene	ND	ND	ND	ND	ND	ND	ND	ND
Dimethylphenylmethylamine	ND	ND	ND	ND	ND	ND	ND	ND
N-Nitroso-di-n-butylamine	ND	ND	ND	ND	ND	ND	ND	ND
4-Chloro-3-methylphenol	ND	ND	ND	ND	ND	ND	ND	ND
Safrole	ND	ND	ND	ND	ND	ND	ND	ND
2-Methylnaphthalene	ND	ND	ND	ND	ND	ND	ND	ND
1,2,4,5-Tetrachlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
Hexachlorocyclopentadiene	ND	ND	ND	ND	ND	ND	ND	ND
2,4,6-Trichlorophenol	ND	ND	ND	ND	ND	ND	ND	ND
Icosafole	ND	ND	ND	ND	ND	ND	ND	ND
2-Chloronaphthalene	ND	ND	ND	ND	ND	ND	ND	ND
2-Nitroaniline	ND	ND	ND	ND	ND	ND	ND	ND
1,4-Naphthoquinone	ND	ND	ND	ND	ND	ND	ND	ND
Dimethylphthalate	ND	ND	ND	ND	ND	ND	ND	ND
1,3-Dinitrobenzene	ND	ND	ND	ND	ND	ND	ND	ND
2,6-Dinitrotoluene	ND	ND	ND	ND	ND	ND	ND	ND
Acenaphthylene	ND	ND	ND	ND	ND	ND	ND	ND
3-Nitroaniline	ND	ND	ND	ND	ND	ND	ND	ND
4-Nitrophenol	ND	ND	ND	ND	ND	ND	ND	ND
2,4-Dinitrophenol	ND	ND	ND	ND	ND	ND	ND	ND
Acenaphthene	ND	ND	ND	ND	ND	ND	ND	ND
2,4-Dinitrotoluene	ND	ND	ND	ND	ND	ND	ND	ND
Dibenzofuran	ND	ND	ND	ND	ND	ND	ND	ND
Pentachlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND
1-Naphthylamine	ND	ND	ND	ND	ND	ND	ND	ND
2-Naphthylamine	ND	ND	ND	ND	ND	ND	ND	ND
2,3,4,6-Tetrachlorophenol	ND	ND	ND	ND	ND	ND	ND	ND
Diethylphthalate	7.303E-04	5.896E-04	1.153E-06	3.345E-07	1.517E-04	7.177E-10	2.370E-06	2.370E-06
4-Chlorophenyl ether	ND	ND	ND	ND	ND	ND	ND	ND

Table B-3: Air Modeling Output Data for Semi-Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Item (grams/m ³) CONC	Pollutant Concentration 1 Item (grams/m ³)	Pollutant Emission Rate (g/sec)/Item	Event Pollutant Emission Rate 1 Item (g/sec)
				M		ER ₁		ER _{EV}	
Fluorene	ND	ND	ND	ND	ND	ND	ND	ND	ND
5-Nitro-o-toluidine	ND	ND	ND	ND	ND	ND	ND	ND	ND
4-Nitroaniline	ND	ND	ND	ND	ND	ND	ND	ND	ND
4,6-Dinitro-2-methylphenol	ND	ND	ND	ND	ND	ND	ND	ND	ND
Diphenylamine/N-NitrosodPDA	ND	ND	ND	ND	ND	ND	ND	ND	ND
Sym-Trinitrobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Diallate	ND	ND	ND	ND	ND	ND	ND	ND	ND
Phenacetin	ND	ND	ND	ND	ND	ND	ND	ND	ND
4-Bromophenylphenyl ether	ND	ND	ND	ND	ND	ND	ND	ND	ND
Hexachlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND
4-Aminobiphenyl	ND	ND	ND	ND	ND	ND	ND	ND	ND
Pronamide	ND	ND	ND	ND	ND	ND	ND	ND	ND
Pentachlorophenol	ND	ND	ND	ND	ND	ND	ND	ND	ND
Pentachloronitrobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Phenanthrene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Anthracene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Carbazole	ND	ND	ND	ND	ND	ND	ND	ND	ND
Di-n-butylphthalate	1.494E-03	1.957E-03	ND	ND	ND	ND	ND	ND	ND
4-Nitroquinoline-1-oxide	ND	ND	ND	ND	ND	ND	ND	ND	ND
Methaprylene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Fluoranthene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Benzidine	ND	ND	ND	ND	ND	ND	ND	ND	ND
Pyrene	ND	ND	ND	ND	ND	ND	ND	ND	ND
p-Dimethylaminoazobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chlorobenzilate	ND	ND	ND	ND	ND	ND	ND	ND	ND
Kepone	ND	ND	ND	ND	ND	ND	ND	ND	ND
Butylbenzylphthalate	3.452E-04	ND	2.829E-06	8.205E-07	3.722E-04	1.761E-09	5.815E-06	5.815E-06	5.815E-06
3,3'-Dimethylbenzidine	ND	ND	ND	ND	ND	ND	ND	ND	ND
2-Acetylaminofluorene	ND	ND	ND	ND	ND	ND	ND	ND	ND
bis(2-Ethylhexyl)phthalate	7.511E-04	5.277E-04	1.831E-06	5.311E-07	2.409E-04	1.140E-09	3.764E-06	3.764E-06	3.764E-06
3,3'-Dichlorobenzidine	ND	ND	ND	ND	ND	ND	ND	ND	ND
Benz(a)anthracene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Chrysene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Di-n-octylphthalate	2.065E-04	ND	1.693E-06	4.909E-07	2.227E-04	1.054E-09	3.479E-06	3.479E-06	3.479E-06
7,12-Dimethylbenz(a)anthracene	ND	ND	ND	ND	ND	ND	ND	ND	ND
Benzo(b)fluoranthene	ND	ND	ND	ND	ND	ND	ND	ND	ND

Table B-3: Air Modeling Output Data for Semi-Volatile Organic Compounds

Compound	Measured Actual Concentration (mg/m ³)	Measured Background Concentration (mg/m ³)	Average Adjusted Emission Factor (lb/lb NEW)	Average Adjusted Emission Factor (lb/item)	Total Mass of Pollutant Emitted (grams/item)	Pollutant Concentration 1 Item (grams/m ³)	Pollutant Emission Rate (g/sec)/item	Event Pollutant Emission Rate 1 Item (g/sec)
				M	M	CONC	ER ₁	ER _{EV}
Benzofluoranthene	ND	ND	ND	ND	ND	ND	ND	ND
Benz(a)pyrene	ND	ND	ND	ND	ND	ND	ND	ND
3-Methylcholanthrene	ND	ND	ND	ND	ND	ND	ND	ND
Indeno(1,2,3-cd)pyrene	ND	ND	ND	ND	ND	ND	ND	ND
Dibenz(a,h)anthracene	ND	ND	ND	ND	ND	ND	ND	ND
Benzo(g,h,i)perylene	ND	ND	ND	ND	ND	ND	ND	ND

Footnotes:

ND = Not Detected

NEW = Net Explosive Weight

APPENDIX C

HEALTH-BASED SCREENING LEVELS AND ACUTE TOXICITY VALUES

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSL)					For the Acute Evaluation (ATV)				
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Region 3 Toxicity Endpoint ($\mu\text{g}/\text{m}^3$)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Value ($\mu\text{g}/\text{m}^3$)	Acute Toxicity	
TSP	12789-66-1	5.00E+01	NA	5.00E+01	NA	NA	NA	NA	NA		
PM ₁₀		5.00E+01	NA	5.00E+01	NA	NA	NA	NA	NA		
HCl	7647-01-0	2.08E+01	nc	2.08E+01	nc	2.08E+01	NA	7.14E+03	T	7.14E+03	
Cl ₂	7782-50-5	2.09E+01	nc	3.65E+02	nc	2.09E+01	2.89E+03	2.90E+03	E	2.89E+03	
Dioxin TEQ	1746-01-6	4.48E-08	c	4.48E-08	c	4.48E-08	NA	3.50E+00	T	3.50E+00	
Carbon Monoxide (CO)	630-08-0	1.57E+02	NA	NA	NA	1.57E+02	2.30E+05	2.28E+05	E	2.30E+05	
Nitrogen Oxide (NOx)	10024-97-2	1.00E+02	NA	NA	NA	1.00E+02	NA	2.70E+05	T	2.70E+05	
HCl (CEM System)	7647-01-0	2.08E+01	nc	2.08E+01	nc	2.08E+01	NA	7.14E+03	T	7.14E+03	
Carbon Dioxide (CO ₂)	124-38-9	NA	NA	NA	NA	NA	NA	5.40E+07	T	5.40E+07	
Sulfur Dioxide (SO ₂)	202-58-84	8.00E+01	NA	8.00E+01	NA	7.89E+02	7.86E+02	E	7.89E+02		
Aluminum	7429-90-5	NA	3.65E+00	nc	3.65E+00	NA	3.00E+04	T	3.00E+04		
Antimony	7440-36-0	NA	1.46E+00	nc	1.46E+00	NA	1.50E+03	T	1.50E+03		
Arsenic	7440-38-2	4.47E-04	c	4.15E-04	c	4.47E-04	NA	3.00E+01	T	3.00E+01	
Barium	7440-39-3	5.21E-01	nc	5.11E-01	nc	5.21E-01	NA	1.50E+03	T	1.50E+03	
Beryllium	7440-41-7	8.00E-04	c	7.45E-04	c	8.00E-04	NA	5.00E+00	T	5.00E+00	
Cadmium	7440-43-9	1.07E-03	c	9.94E-04	c	1.07E-03	NA	3.00E+01	T	3.00E+01	
Chromium	7440-43-9	NA	c	1.53E-04	c	1.53E-04	NA	1.50E+03	T	1.50E+03	
Cobalt	7440-48-4	NA		2.20E+02	nc	2.20E+02	NA	6.00E+01	T	6.00E+01	
Copper	7440-50-8	NA		1.46E+02	nc	1.46E+02	NA	3.00E+03	T	3.00E+03	
Lead	7439-92-1	1.50E+00		NA		1.50E+00	NA	1.50E+02	T	1.50E+02	
Magnesium	7439-95-4	NA		NA		NA	NA	3.00E+04	T	3.00E+04	
Manganese	7439-96-5	5.11E-02	nc	5.22E-02	nc	5.11E-02	NA	3.00E+03	T	3.00E+03	
Nickel	7440-02-0	NA		7.30E+01	nc	7.30E+01	NA	3.00E+03	T	3.00E+03	
Phosphorus	7723-14-0	NA		NA		NA	NA	3.00E+02	T	3.00E+02	
Selenium	7782-49-2	NA		1.83E+01	nc	1.83E+01	NA	6.00E+02	T	6.00E+02	
Silver	7740-22-4	NA		1.83E+01	nc	1.83E+01	NA	3.00E+02	T	3.00E+02	
Thallium	7740-28-0	NA		2.56E-01	nc	2.56E-01	NA	3.00E+02	T	3.00E+02	
Zinc	7740-66-6	NA		1.10E+03	nc	1.10E+03	NA	3.00E+04	T	3.00E+04	
Mercury	7439-97-6	3.13E-01	nc	3.14E-01	nc	3.13E-01	NA	1.00E+02	T	1.00E+02	
TNMHC		NA		NA		NA	NA	NA	NA		
Ethane	74-84-0	NA		NA		NA	NA	NA	NA		
Ethylene	74-85-1	NA		NA		NA	NA	4.60E+05	T	4.60E+05	
Acetylene	74-86-2	NA		NA		NA	NA	NA	NA		
Propane	74-98-6	NA		NA		NA	NA	3.78E+06	T	3.78E+06	
Propene	115-07-1	NA		NA		NA	NA	NA	NA		
i-Butane	106-97-8	NA		NA		NA	NA	5.71E+06	T	5.71E+06	
i-Butene	25167-67-3	NA		NA		NA	NA	NA	NA		
1-Butene	106-98-9	NA		NA		NA	NA	NA	NA		
1,3-Butadiene	106-99-0	3.74E-03	c	3.48E-03	c	3.74E-03	NA	2.20E+04	2.21E+04	E	2.20E+04

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSL)						For the Acute Evaluation (ATV)		
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)
n-Butane	106-97-8	NA	NA	NA	NA	NA	NA	NA	5.71E+06	5.71E+06
trans-2-Butene	624-64-6	NA	NA	NA	NA	NA	NA	NA	NA	NA
2,2-Dimethylpropane	463-82-1	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-2-Butene	590-18-1	NA	NA	NA	NA	NA	NA	NA	NA	NA
3-Methyl-1-butene	563-45-1	NA	NA	NA	NA	NA	NA	NA	NA	NA
i-Pentane	109-66-0	NA	NA	NA	NA	NA	NA	NA	1.80E+06	T
1-Pentene	109-67-1	NA	NA	NA	NA	NA	NA	NA	NA	1.80E+06
2-Methyl-1-butene	563-46-2	NA	NA	NA	NA	NA	NA	NA	NA	NA
n-Pentane	109-66-0	NA	NA	NA	NA	NA	NA	NA	1.80E+06	T
Isoprene	78-79-5	NA	NA	NA	NA	NA	NA	NA	NA	NA
trans-2-Pentene	646-04-8	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-2-Pentene	627-20-3	NA	NA	NA	NA	NA	NA	NA	NA	NA
2-Methyl-2-butene	513-35-9	NA	NA	NA	NA	NA	NA	NA	NA	NA
2,2-Dimethylbutane	75-83-2	NA	NA	NA	NA	NA	NA	NA	1.80E+06	T
Cyclopentene	142-29-0	NA	NA	NA	NA	NA	NA	NA	NA	NA
4-Methyl-1-pentene	691-37-2	NA	NA	NA	NA	NA	NA	NA	NA	NA
Cyclopentane	287-92-3	NA	NA	NA	NA	NA	NA	NA	NA	NA
2,3-Dimethylbutane	79-29-8	NA	NA	NA	NA	NA	NA	NA	1.80E+06	T
cis-4-Methyl-2-pentene	691-38-3	NA	NA	NA	NA	NA	NA	NA	NA	NA
2-Methylpentane	107-83-5	NA	NA	NA	NA	NA	NA	NA	NA	NA
3-Methylpentane	96-14-0	NA	NA	NA	NA	NA	NA	NA	NA	NA
2-Methyl-1-pentene	763-29-1	NA	NA	NA	NA	NA	NA	NA	NA	NA
1-Lexene	592-41-6	NA	NA	NA	NA	NA	NA	NA	1.03E+05	T
n-Hexane	110-54-3	2.10E+02	nc	2.10E+02	nc	2.10E+02	nc	2.10E+02	NA	5.28E+05
trans-2-Hexene	4050-45-7	NA	NA	NA	NA	NA	NA	NA	NA	NA
2-Methyl-2-pentene	625-27-4	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-2-Hexene	7688-21-3	NA	NA	NA	NA	NA	NA	NA	NA	NA
Methylcyclopentane	96-37-7	NA	NA	NA	NA	NA	NA	NA	NA	NA
2,4-Dimethylpentane	108-08-7	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	71-43-2	2.50E-01	c	2.2E-01	c	2.50E-01	c	1.56E+05	E	1.56E+05
Cyclohexane	110-82-7	NA	NA	NA	NA	NA	NA	3.10E+06	T	3.10E+06
2-Methylhexane	591-76-4	NA	NA	NA	NA	NA	NA	NA	NA	NA
2,3-Dimethylpentane	565-59-3	NA	NA	NA	NA	NA	NA	NA	NA	NA
3-Methylhexane	589-34-4	NA	NA	NA	NA	NA	NA	NA	NA	NA
2,2,4-Trimethylpentane	540-84-1	NA	NA	NA	NA	NA	NA	3.50E+05	T	3.50E+05
n-Heptane	142-82-5	NA	NA	NA	NA	NA	NA	1.80E+06	T	1.80E+06
2,4,4-Trimethyl-1-pentene	107-39-1	NA	NA	NA	NA	NA	NA	NA	NA	NA
Methylcyclohexane	108-87-2	3.10E+03	nc	3.10E+03	nc	3.10E+03	nc	NA	4.81E+06	T
2,4,4-Trimethyl-2-pentene	107-40-4	NA	NA	NA	NA	NA	NA	NA	NA	NA

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSEL)				For the Acute Evaluation (ATV)			
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (C or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (C or nc)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)
2,5-Dimethylhexane	592-13-2	NA	NA	NA	NA	NA	NA	NA	NA
2,4-Dimethylhexane	569-43-5	NA	NA	NA	NA	NA	NA	NA	NA
2,3,4-Trimethylpentane	565-59-3	NA	NA	NA	NA	NA	NA	NA	NA
Toluene	108-88-3	4.02E+02	nc	4.16E+02	nc	4.02E+02	1.88E+05	1.89E+05	E
2,3-Dimethylhexane	564-94-1	NA	NA	NA	NA	NA	NA	NA	NA
2-Methylheptane	592-27-8	NA	NA	NA	NA	NA	NA	NA	NA
3-Ethylhexane	619-99-8	NA	NA	NA	NA	NA	NA	NA	NA
2,2-Dimethylheptane	1071-26-7	NA	NA	NA	NA	NA	NA	NA	NA
2,2,4-Trimethylhexane	16747-26-5	NA	NA	NA	NA	NA	NA	NA	NA
n-Octane	111-65-9	NA	NA	NA	NA	NA	NA	NA	NA
Ethylcyclohexane	1678-91-7	NA	NA	NA	NA	NA	NA	NA	NA
Ethylbenzene	100-41-4	1.10E+03	nc	1.1E+03	nc	1.10E+03	NA	5.43E+05	T
m-Xylene & p-Xylene	108-38-3	NA	NA	NA	NA	NA	NA	6.51E+05	T
Styrene	100-42-5	1.10E+03	nc	1.0E+03	nc	1.10E+03	2.13E+05	2.13E+05	E
o-Xylene	95-47-6	NA	NA	NA	NA	7.30E+03	NA	6.51E+05	T
n-Nonane	111-84-2	NA	NA	NA	NA	4.0E+02	NA	4.02E+02	T
i-Propylbenzene	98-82-8	4.00E+02	nc	4.0E+02	nc	4.00E+02	NA	4.00E+02	NA
n-Propylbenzene	103-65-1	3.65E+01	nc	1.5E+02	nc	3.65E+01	NA	NA	NA
p-Ethyltoluene	622-96-8	NA	NA	NA	NA	NA	NA	1.25E+05	T
m-Ethyltoluene	620-14-4	NA	NA	NA	NA	NA	NA	NA	NA
1,3,5-Trimethylbenzene	108-67-8	6.20E+00	nc	6.2E+00	nc	6.20E+00	NA	3.68E+05	T
o-Ethyltoluene	611-14-3	NA	NA	NA	NA	NA	NA	7.50E+02	T
1,2,4-Trimethylbenzene & sec-Butylbenzene	95-63-6	6.21E+00	nc	6.21E+00	nc	6.21E+00	NA	1.80E+05	T
n-Decane	124-18-5	NA	NA	NA	NA	NA	NA	4.37E+03	T
alpha-Pinene	80-56-8	NA	NA	NA	NA	NA	NA	4.00E+04	T
beta-Pinene	127-91-3	NA	NA	NA	NA	NA	NA	NA	NA
delta-3-Carene	13466-78-9	NA	NA	NA	NA	NA	NA	NA	NA
d-Limonene	5989-27-5	NA	NA	NA	NA	NA	NA	1.95E+06	T
MTBE	1634-04-4	3.10E+03	nc	3.1E+03	nc	3.10E+03	NA	4.32E+05	T
Dichlorodifluoromethane	75-71-8	2.10E+02	nc	1.8E+02	nc	2.10E+02	NA	1.48E+07	T
Methylchloride	74-87-33	NA	NA	NA	NA	NA	NA	NA	NA
Dichlorotetrafluoroethane	374-07-2	NA	NA	NA	NA	NA	NA	NA	NA
Chloroethene	75-01-4	2.20E-02	C	2.1E-02	C	2.20E-02	NA	1.28E+04	T
1,3-Butadiene	106-99-0	3.74E-03	C	3.48E-03	C	3.74E-03	2.20E+04	2.21E+04	E
Methylbromide	74-83-9	5.20E+00	nc	5.1E+00	nc	5.20E+00	NA	5.82E+04	T
Ethylchloride	75-00-3	2.30E+00	C	2.2E+00	C	2.30E+00	NA	7.92E+06	T
Trichloromonofluoromethane	75-69-4	7.30E+02	nc	7.30E+02	nc	7.30E+02	NA	2.81E+06	T
Vinyldiene chloride	75-35-4	NA	NA	NA	NA	NA	NA	7.92E+04	T
Methylene chloride	75-09-2	4.10E+00	C	3.8E+00	C	4.10E+00	6.94E+05	6.96E+05	E

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	PRG ($\mu\text{g}/\text{m}^3$)	For the Chronic Evaluation (HBSL)				For the Acute Evaluation (ATV)			
			Region 9 Toxicity Endpoint (c or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)
Allyl chloride	107-05-1	1.00E+00	nc	NA	1.00E+00	9.39E+03	9.39E+03	E	9.39E+03	
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	3.13E+04	nc	3.14E+04	3.13E+04	NA	9.58E+06	T	9.58E+06	
1,1-Dichloroethane	75-34-3	5.21E+02	nc	5.11E+02	5.21E+02	NA	1.21E+06	T	1.21E+06	
1,2-Dichloroethene	540-59-0	NA	3.29E+01	nc	3.29E+01	NA	2.38E+06	T	2.38E+06	
Chloroform	67-66-3	8.40E-02	c	2.2E+00	c	8.40E-02	NA	9.76E+03	T	9.76E+03
1,2-Dichloroethane	107-06-2	7.39E-02	c	6.88E-02	c	7.39E-02	NA	8.08E+03	T	8.08E+03
Methylchloroform	71-55-6	1.00E+03	nc	2.3E+03	nc	1.00E+03	NA	1.91E+06	T	1.91E+06
Benzene	71-43-2	2.50E-01	c	2.2E-01	c	2.50E-01	NA	1.60E+05	T	1.60E+05
Carbon tetrachloride	56-23-5	1.04E+03	nc	1.04E+03	nc	1.04E+03	1.28E+05	1.26E+05	E	1.28E+05
1,2-Dichloropropane	78-87-5	9.89E-02	c	9.21E-02	c	9.89E-02	NA	5.08E+05	T	5.08E+05
Trichloroethylene	79-01-6	1.12E+00	c	1.04E+00	c	1.12E+00	NA	5.37E+05	T	5.37E+05
cis 1,3-Dichloro-1-propene	10061-01-5	NA	NA	NA	NA	NA	1.14E+04	T	1.14E+04	
trans 1,3-Dichloro-1-propene	10061-02-6	NA	NA	NA	NA	NA	NA	NA	NA	
1,1,2-Trichloroethane	79-00-5	1.20E-01	c	1.12E-01	c	1.20E-01	NA	1.64E+05	T	1.64E+05
Toluene	108-88-3	4.02E+02	nc	4.16E+02	nc	4.02E+02	1.88E+05	1.89E+05	E	1.88E+05
1,2-Dibromoethane	106-93-4	8.73E-03	c	8.24E-03	c	8.73E-03	NA	1.54E+05	T	1.54E+05
Perchloroethylene	127-18-4	3.31E+00	c	3.13E+00	c	3.31E+00	6.89E+05	6.78E+05	E	6.89E+05
Chlorobenzene	108-90-7	6.20E+01	nc	6.2E+01	nc	6.20E+01	NA	1.38E+05	T	1.38E+05
Ethylbenzene	100-41-4	1.10E+03	nc	1.1E+03	nc	1.10E+03	NA	4.34E+03	T	4.34E+03
m&p-Xylene	108-38-3	7.30E+02	nc	NA	NA	7.30E+02	NA	6.51E+05	T	6.51E+05
Styrene	100-42-5	1.06E+03	nc	1.04E+03	nc	1.06E+03	2.13E+05	2.13E+05	E	2.13E+05
1,1,2,2-Tetrachloroethane	79-34-5	3.31E-02	c	3.13E-02	c	3.31E-02	NA	2.06E+04	T	2.06E+04
o-Xylene	95-47-6	7.30E+02	nc	7.30E+03	nc	7.30E+02	NA	6.51E+05	T	6.51E+05
p-Ethyltoluene	622-96-8	NA	NA	NA	NA	NA	1.25E+05	T	1.25E+05	
1,3,5-Trimethylbenzene	108-67-8	6.21E+00	nc	6.21E+00	nc	6.21E+00	NA	3.68E+05	T	3.68E+05
1,2,4-Trimethylbenzene	95-63-6	6.21E+00	nc	6.21E+00	nc	6.21E+00	NA	1.80E+05	T	1.80E+05
Benzyl chloride	100-44-7	4.00E-02	nc	3.7E-02	c	4.00E-02	5.20E+03	5.17E+03	E	5.20E+03
m-Dichlorobenzene	541-73-1	3.30E+00	nc	3.3E+00	nc	3.30E+00	NA	3.61E+04	T	3.61E+04
p-Dichlorobenzene	106-46-7	2.80E-01	c	2.85E-01	c	2.80E-01	NA	6.61E+05	T	6.61E+05
o-Dichlorobenzene	95-50-1	2.09E+02	nc	3.29E+01	nc	2.09E+02	NA	3.01E+05	T	3.01E+05
1,2,4-Trichlorobenzene	120-82-1	NA	NA	NA	NA	NA	3.71E+04	T	3.71E+04	
Hexachlorobutadiene	87-68-3	8.73E-02	c	8.03E-02	c	8.73E-02	3.21E+04	3.20E+04	E	3.21E+04
trans-1,2-Dichloroethylene	156-60-5	7.30E+01	nc	7.30E+01	nc	7.30E+01	NA	4.95E+04	T	4.95E+04
o-Chlorotoluene	95-49-8	7.30E+01	nc	7.30E+01	nc	7.30E+01	NA	3.88E+05	T	3.88E+05
p-Chlorotoluene	106-43-4	NA	NA	NA	NA	NA	3.88E+05	T	3.88E+05	
1,3,5-Trichlorobenzene	108-70-3	NA	NA	NA	NA	NA	NA	NA	NA	
1,2,3-Trichlorobenzene	87-61-6	NA	NA	NA	NA	NA	5.00E+04	T	5.00E+04	
Methyl nitrite	624-91-9	NA	NA	NA	NA	NA	NA	NA	NA	
Acetonitrile	75-05-8	6.20E+01	nc	6.2E+01	nc	6.20E+01	NA	1.01E+05	T	1.01E+05

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSL)						For the Acute Evaluation (ATV)					
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Value ($\mu\text{g}/\text{m}^3$)	Acute Toxicity		
Acrylonitrile	107-13-1	2.80E-02	C	2.6E-02	C	2.80E-02	2.20E+04	2.17E+04	E	2.20E+04			
Nitromethane	75-52-5	NA	NA	NA	NA	NA	NA	1.50E+05	T	1.50E+05			
Benzonitrile	100-47-0	NA	NA	NA	NA	NA	NA	1.50E+04	T	1.50E+04			
Nitrobenzene	98-95-3	2.09E+00	nc	2.19E+00	nc	2.09E+00	NA	NA	1.51E+04	T	1.51E+04		
Carbonyl Sulfide	463-58-1	NA	NA	NA	NA	NA	NA	9.84E+03	T	9.84E+03			
Sulfur Dioxide	7446-09-5	NA	NA	NA	NA	NA	7.80E+02	7.86E+02	E	7.80E+02			
Carbon Disulfide	75-15-0	7.30E+02	nc	7.3E+02	nc	7.30E+02	NA	3.73E+04	T	3.73E+04			
Thiophene	110-02-1	NA	NA	NA	NA	NA	NA	NA	NA	NA			
Dimethyl Disulfide	624-92-0	NA	NA	NA	NA	NA	4.00E+01	3.85E+01	E	4.00E+01			
2-Methyl Thiophene	554-14-3	NA	NA	NA	NA	NA	NA	NA	NA	NA			
3-Methyl Thiophene	616-44-4	NA	NA	NA	NA	NA	NA	NA	NA	NA			
Dimethyl Trisulfide	3658-80-8	NA	NA	NA	NA	NA	NA	NA	NA	NA			
Isothiocyanomethane	556-61-6	NA	NA	NA	NA	NA	NA	NA	NA	NA			
2-Chlorothiophene	96-43-5	NA	NA	NA	NA	NA	NA	NA	NA	NA			
3-Chlorothiophene	17249-80-8	NA	NA	NA	NA	NA	NA	NA	NA	NA			
2-Thiophenecarboxaldehyde	98-03-3	NA	NA	NA	NA	NA	NA	NA	NA	NA			
Naphthalene	91-20-3	3.13E+00	nc	3.29E+00	nc	3.13E+00	NA	7.86E+04	T	7.86E+04			
Acetaldehyde	75-07-0	8.70E-01	C	8.1E-01	C	8.70E-01	1.80E+04	1.80E+04	E	1.80E+04			
Acrolein	107-02-8	2.10E-02	nc	2.1E-02	nc	2.10E-02	2.30E+02	2.29E+03	E	2.30E+02			
Acetone	67-64-1	3.40E+02	nc	3.7E+02	nc	3.40E+02	NA	2.37E+06	T	2.37E+06			
Propanal	123-38-6	NA	NA	NA	NA	NA	NA	7.50E+04	T	7.50E+04			
Furan	110-00-9	3.70E+00	nc	NA	NA	3.70E+00	NA	1.67E+02	T	1.67E+02			
2-Propanol	67-63-0	NA	NA	NA	NA	NA	NA	9.84E+05	T	9.84E+05			
2-Methylpropanal	78-84-2	NA	NA	NA	NA	NA	NA	NA	NA	NA			
Methacrolein	78-85-3	NA	NA	NA	NA	NA	NA	NA	NA	NA			
2,3-Butanedione	625-34-3	NA	NA	NA	NA	NA	NA	NA	NA	NA			
Methyl-Vinyl Ketone	78-94-4	NA	NA	NA	NA	NA	NA	NA	NA	NA			
MTBE	1634-04-4	3.10E+03	nc	3.1E+03	nc	3.10E+03	NA	4.32E+05	T	4.32E+05			
Butanal	123-72-8	NA	NA	NA	NA	NA	NA	7.38E+04	T	7.38E+04			
2-Butanone	78-93-3	1.00E+03	nc	1.0E+03	nc	1.00E+03	NA	8.85E+05	T	8.85E+05			
Tetrahydrofuran	109-99-9	9.89E-01	nc	9.21E-01	C	9.89E-01	NA	7.38E+05	T	7.38E+05			
2-Methyl-1-propanol	78-83-1	1.10E+03	nc	1.1E+03	nc	1.10E+03	NA	4.55E+05	T	4.55E+05			
trans-2-Butenal	123-73-9	3.54E-03	C	3.30E-03	C	3.54E-03	NA	NA	NA	NA			
Acetic Acid	64-19-7	NA	NA	NA	NA	NA	NA	3.68E+04	T	3.68E+04			
2-Pentanone	107-87-9	NA	NA	NA	NA	NA	NA	8.80E+05	T	8.80E+05			
Pentanal	110-62-3	NA	NA	NA	NA	NA	NA	NA	NA	NA			
4-Methyl-2-pentanone	108-10-1	8.30E+01	nc	7.3E+01	nc	8.30E+01	NA	3.07E+05	T	3.07E+05			
trans-2-Pentenal	1567-87-0	NA	NA	NA	NA	NA	NA	NA	NA	NA			
Cyclopentanone	120-92-3	NA	NA	NA	NA	NA	NA	NA	NA	NA			

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSL)					For the Acute Evaluation (ATV)				
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	
2-Hexanone	591-78-6	NA	5.1E+00	nc	5.11E+00	NA	4.09E+04	NA	T	4.09E+04	
Hexanal	66-25-1	NA	NA	NA	NA	NA	NA	NA	NA	NA	
3-Furaidehyde	498-60-2	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Butyl/Acetate	123-86-4	NA	NA	NA	NA	NA	NA	NA	NA	NA	
2-Furaidehyde	98-01-1	5.20E+01	nc	3.7E+01	nc	5.20E+01	NA	7.86E+03	T	7.86E+03	
trans-2-Hexenal	6728-26-3	NA	NA	NA	NA	NA	NA	NA	NA	NA	
1-Hexanol	111-27-3	NA	NA	NA	NA	NA	NA	8.36E+03	T	8.36E+03	
3-Heptanone	106-35-4	NA	NA	NA	NA	NA	NA	NA	NA	NA	
2-Heptanone	110-43-0	NA	NA	NA	NA	NA	NA	1.70E+03	T	1.70E+03	
Heptanal	66-25-1	NA	NA	NA	NA	NA	NA	NA	NA	NA	
trans-2-Heptenal	18829-55-5	NA	NA	NA	NA	NA	NA	NA	NA	NA	
5-Methyl-2-furaldehyde	620-02-0	NA	NA	NA	NA	NA	NA	NA	NA	NA	
6-Methyl-2-heptanone	928-68-7	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Benzaldehyde	100-52-7	3.70E+02	nc	3.7E+02	nc	3.70E+02	NA	1.50E+04	T	1.50E+04	
1-Heptanol	111-70-6	NA	NA	NA	NA	NA	NA	NA	NA	NA	
6-Methyl-5-hepten-2-one	110-93-0	NA	NA	NA	NA	NA	NA	NA	NA	NA	
2-Octanone	111-13-7	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Octanal	124-13-0	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Benzofuran	271-89-6	NA	NA	NA	NA	NA	NA	NA	NA	NA	
trans-2-Octenol	254-8-87-0	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Acetophenone	98-86-2	2.10E-02	nc	2.1E-02	nc	2.10E-02	NA	3.00E+04	T	3.00E+04	
2-Nonanone	821-55-6	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Nonanal	124-19-6	NA	NA	NA	NA	NA	NA	NA	NA	NA	
trans-2-Nonenal	18829-56-6	NA	NA	NA	NA	NA	NA	NA	NA	NA	
2-Decanone	693-54-9	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Decanal	112-31-2	NA	NA	NA	NA	NA	NA	NA	NA	NA	
N-Nitrosodimethylamine	62-75-9	1.40E-04	c	1.2E-04	c	1.40E-04	NA	2.50E+03	T	2.50E+03	
Pyridine	110-86-1	3.65E+00	nc	3.65E+00	nc	3.65E+00	NA	4.85E+04	T	4.85E+04	
2-Picoline	109-06-8	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Methyl methanesulfonate	66-27-3	NA	NA	NA	NA	NA	NA	NA	NA	NA	
N-Nitrosomethylamine	10595-95-6	3.06E-04	c	2.85E-04	c	3.06E-04	NA	NA	NA	NA	
N-Nitrosodiethylamine	55-18-5	4.47E-05	c	4.17E-05	c	4.47E-05	NA	NA	NA	NA	
Ethyl methanesulfonate	62-50-0	NA	NA	NA	NA	NA	NA	NA	NA	NA	
Phenol	108-95-2	2.19E+03	nc	2.19E+03	nc	2.19E+03	NA	3.85E+04	E	3.85E+04	
Aniline	62-53-3	NA	1.1E+00	nc	1.06E+00	NA	2.29E+04	T	2.29E+04		
bis(2-Chloroethyl)ether	111-44-4	5.80E-03	c	5.7E-03	c	5.80E-03	NA	5.85E+04	T	5.85E+04	
Pentachloroethane	76-01-7	NA	NA	NA	NA	NA	NA	NA	NA	NA	
2-Chlorophenol	95-57-8	1.80E+01	nc	1.8E+01	nc	1.80E+01	NA	5.25E+03	T	5.25E+03	
1,3-Dichlorobenzene	543-73-1	NA	NA	NA	NA	NA	NA	NA	NA	NA	

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSL)					For the Acute Evaluation (ATV)				
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	
1,4-Dichlorobenzene	106-46-7	2.80E-01	c	2.85E-01	c	2.80E-01	NA	6.61E+05	T	6.61E+05	
Benzyl alcohol	100-51-6	1.10E+03	nc	1.1E+03	nc	1.10E+03	NA	5.53E+04	T	5.53E+04	
2-Methylphenol	95-48-7	NA	NA	NA	NA	NA	NA	6.63E+04	T	6.63E+04	
1,2-Dichlorobenzene	95-50-1	2.09E+02	nc	3.29E+01	nc	2.09E+02	NA	3.01E+05	T	3.01E+05	
bis(2-Chloroisopropyl)ether	108-60-1	1.92E-01	c	1.79E-01	c	1.92E-01	NA	6.99E+04	T	6.99E+04	
o-Tolidine	95-53-4	2.80E-02	c	2.6E-02	c	2.80E-02	NA	2.63E+04	T	2.63E+04	
4-Methylphenol/3-Methylphenol	1319-77-3	NA	NA	NA	NA	NA	NA	6.63E+04	T	6.63E+04	
N-Nitroso-di-n-propylamine	621-64-7	9.61E-04	c	8.94E-04	c	9.61E-04	NA	5.32E+03	T	5.32E+03	
Acetophenone	98-86-2	2.10E-02	nc	2.1E-02	nc	2.10E-02	NA	1.47E+05	T	1.47E+05	
N-Nitrosomorpholine	59-89-2	NA	NA	NA	NA	NA	NA	3.00E+04	T	3.00E+04	
N-Nitrosopyrrolidine	930-55-2	3.15E-03	c	3.0E-03	c	3.15E-03	NA	NA			
Hexachloroethane	67-72-1	4.80E-01	c	4.47E-01	c	4.80E-01	NA	2.90E+04	T	2.90E+04	
Nitrobenzene	98-95-3	2.09E+00	nc	2.19E+00	nc	2.09E+00	NA	1.51E+04	T	1.51E+04	
N-Nitrosopiperidine	100-75-4	NA	NA	NA	NA	NA	NA				
Isophorone	78-59-1	7.08E+00	c	6.59E+00	c	7.08E+00	NA	2.83E+04	T	2.83E+04	
2,4-Dimethylphenol	105-67-9	7.30E+01	nc	7.3E+01	nc	7.30E+01	NA	NA			
2-Nitrophenol	88-75-5	NA	NA	NA	NA	NA	NA				
bis(2-Chloroethoxy)methane	111-91-1	NA	NA	NA	NA	NA	NA				
Benzoic acid	65-85-0	1.50E+04	nc	1.5E+04	nc	1.50E+04	NA	1.25E+04	T	1.25E+04	
2,4-Dichlorophenol	120-83-2	1.10E+01	nc	1.1E+01	nc	1.10E+01	NA	3.00E+04	T	3.00E+04	
1,2,4-Trichlorobenzene	120-82-1	NA	NA	NA	NA	NA	NA	3.71E+04	T	3.71E+04	
Naphthalene	91-20-3	3.13E+00	nc	3.29E+00	nc	3.13E+00	NA	7.86E+04	T	7.86E+04	
p-Chloroaniline	106-47-8	1.46E+01	nc	1.46E+01	nc	1.46E+01	NA	5.21E+03	T	5.21E+03	
2,6-Dichlorophenol	87-65-0	NA	NA	NA	NA	NA	NA	3.00E+04	T	3.00E+04	
Hexachloropropene	1888-71-7	NA	NA	NA	NA	NA	NA				
Hexachlorobutadiene	87-68-3	8.73E-02	c	8.03E-02	c	8.73E-02	3.21E+04	3.20E+04	E	3.21E+04	
Dimethylphenylethylamine	NA	NA	NA	NA	NA	NA	NA				
N-Nitroso-di-n-butylamine	924-16-3	1.20E-03	c	1.12E-03	c	1.20E-03	NA	2.00E+04	T	2.00E+04	
4-Chloro-3-methylphenol	35421-08-0	NA	NA	NA	NA	NA	NA	3.00E+04	T	3.00E+04	
Safrole	94-59-7	NA	NA	NA	NA	NA	NA				
2-Methylnaphthalene	91-57-6	NA	NA	NA	NA	NA	NA				
1,2,4,5-Tetrachlorobenzene	95-94-3	1.10E+00	nc	1.10E+00	nc	1.10E+00	NA	3.00E+04	T	3.00E+04	
Hexachlorocyclopentadiene	77-47-4	7.30E-02	nc	7.30E-02	nc	7.30E-02	NA	2.23E+02	T	2.23E+02	
2,4,6-Trichlorophenol	88-06-2	6.20E-01	c	6.3E-01	c	6.20E-01	NA	3.00E+04	T	3.00E+04	
2,4,5-Trichlorophenol	95-95-4	3.70E+02	nc	3.7E+02	nc	3.70E+02	NA	3.00E+04	T	3.00E+04	
Iosafrole	120-58-1	NA	NA	NA	NA	NA	NA				
2-Chloronaphthalene	91-58-7	2.90E-02	nc	2.9E+02	nc	2.90E+02	NA	6.00E+02	T	6.00E+02	
2-Nitroaniline	88-74-4	2.10E-01	nc	2.1E-01	nc	2.10E-01	NA	NA			
1,4-Naphthoquinone	130-15-4	NA	NA	NA	NA	NA	NA	2.50E+02	T	2.50E+02	

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSL)				For the Acute Evaluation (ATV)				
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Health-based Screening Level (c or nc) ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	
Dimethylphthalate	131-11-3	3.65E+04	nc	3.65E+04	nc	3.65E+04	NA	1.50E+04	T	1.50E+04
1,3-Dinitrobenzene	99-65-0	3.70E-01	nc	3.7E-01	nc	3.70E-01	NA	3.00E+03	T	3.00E+03
2,6-Dinitrotoluene	606-20-2	3.70E+00	nc	3.7E+00	nc	3.70E+00	NA	6.00E+02	T	6.00E+02
Aceanaphthylene	208-96-8	NA	NA	NA	NA	NA	2.00E+02	T	2.00E+02	
3-Nitroaniline	99-09-2	NA	NA	NA	NA	NA	NA	NA	NA	
4-Nitrophenol	100-02-7	2.90E+01	nc	2.9E+01	nc	2.90E+01	NA	3.00E+04	T	3.00E+04
2,4-Dinitrophenol	51-28-5	7.30E+00	nc	7.3E+00	nc	7.30E+00	NA	7.50E+03	T	7.50E+03
Acenaphthene	83-32-9	2.20E+02	nc	2.2E+02	nc	2.20E+02	NA	1.25E+03	T	1.25E+03
2,4-Dinitrotoluene	121-14-2	7.30E+00	nc	7.3E+00	nc	7.30E+00	NA	6.00E+02	T	6.00E+02
Dibenzofuran	132-64-9	1.46E+01	nc	1.46E+01	nc	1.46E+01	NA	1.50E+00	T	1.50E+00
Pentachlorobenzene	608-93-5	2.92E+00	nc	2.92E+00	nc	2.92E+00	NA	3.00E+04	T	3.00E+04
1-Naphthylamine	134-32-7	NA	NA	NA	NA	NA	NA	3.50E+04	T	3.50E+04
2-Naphthylamine	91-59-8	NA	NA	NA	NA	NA	NA	7.50E+03	T	7.50E+03
2,3,4,6-Tetrachlorophenol	58-90-2	1.10E+02	nc	1.1E+02	nc	1.10E+02	NA	NA	NA	NA
Diethylphthalate	84-66-2	2.92E+03	nc	2.92E+03	nc	2.92E+03	NA	1.50E+04	T	1.50E+04
4-Chlorophenylphenyl ether	7005-72-3	NA	NA	NA	NA	NA	NA	NA	NA	NA
Fluorene	86-73-7	1.46E+02	nc	1.46E+02	nc	1.46E+02	NA	7.50E+04	T	7.50E+04
5-Nitro-o-tolidine	99-55-8	2.00E-01	c	1.9E-01	c	2.00E-01	NA	NA	NA	NA
4-Nitroaniline	100-01-6	NA	NA	NA	NA	NA	NA	9.00E+03	T	9.00E+03
4,6-Dinitro-2-methylphenol	534-52-1	NA	NA	3.7E-01	nc	3.65E-01	NA	5.00E+02	T	5.00E+02
Diphenylamine/N-NitrosoDPA	62-75-9	NA	NA	NA	NA	NA	NA	2.50E+03	T	2.50E+03
sym-Trinitrotoluene	99-35-4	1.10E+02	nc	1.10E+02	nc	1.10E+02	NA	3.00E+04	T	3.00E+04
Diallate	2303-16-4	1.10E-01	c	NA	NA	1.10E-01	NA	NA	NA	NA
Phenacetin	62-44-2	NA	NA	NA	NA	NA	NA	3.00E+04	T	3.00E+04
4-Bromophenylphenyl ether	101-55-3	NA	NA	NA	NA	NA	NA	NA	NA	NA
Hexachlorobenzene	118-74-1	4.18E-03	c	3.91E-03	c	4.18E-03	NA	7.50E+01	T	7.50E+01
4-Aminobiphenyl	92-67-1	NA	NA	NA	NA	NA	NA	1.50E+03	T	1.50E+03
Promamide	23950-58-5	2.74E+02	nc	NA	NA	2.74E+02	NA	NA	NA	NA
Pentachlorophenol	87-86-5	5.60E-02	c	5.22E-02	c	5.60E-02	NA	1.50E+03	T	1.50E+03
Pentachloronitrobenzene	82-68-8	2.59E-02	c	2.41E-02	c	2.59E-02	NA	1.50E+03	T	1.50E+03
Phenanthrene	85-01-8	NA	NA	NA	NA	NA	NA	2.00E+03	T	2.00E+03
Anthracene	120-12-7	1.10E+03	nc	1.1E+03	nc	1.10E+03	NA	6.00E+03	T	6.00E+03
Carbazole	86-74-8	3.36E-01	c	3.13E-01	c	3.36E-01	NA	NA	NA	NA
Di-n-butylphthalate	84-74-2	3.65E+02	nc	3.65E+02	nc	3.65E+02	NA	1.50E+04	T	1.50E+04
4-Nitroquinoline-1-oxide	56-57-5	NA	NA	NA	NA	NA	NA	NA	NA	NA
Methaphylylene	91-80-5	NA	NA	NA	NA	NA	NA	NA	NA	NA
Fluoranthene	206-44-0	1.50E+02	nc	1.5E+02	nc	1.50E+02	NA	3.00E+01	T	3.00E+01
Benzidine	92-87-5	2.90E-05	c	NA	NA	2.90E-05	NA	5.00E+02	T	5.00E+02
Pyrene	129-00-0	NA	NA	NA	NA	NA	NA	1.50E+04	T	1.50E+04

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

Compound	CAS #	For the Chronic Evaluation (HBSL)					For the Acute Evaluation (ATV)				
		Region 9 PRG ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Region 3 RBC ($\mu\text{g}/\text{m}^3$)	Toxicity Endpoint (c or nc)	Health-based Screening Level ($\mu\text{g}/\text{m}^3$)	ERPG ($\mu\text{g}/\text{m}^3$)	TEEL ($\mu\text{g}/\text{m}^3$)	Source (T or E)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	
p-Dimethylaminoazobenzene	60-11-7	NA	NA	2.32E-02	c	2.49E-02	NA	NA	7.50E+04	T	7.50E+04
Chlorobenzilate	510-15-6	2.49E-02	c	NA	NA	2.49E-02	NA	NA	2.50E+02	T	2.50E+02
Kepone	143-50-0	3.74E-04	c	NA	NA	3.74E-04	NA	NA	1.00E+02	T	1.00E+02
Butylbenzylphthalate	85-68-7	7.30E+02	nc	7.30E+02	nc	7.30E+02	NA	NA	5.00E+05	T	5.00E+05
3,3'-Dimethylbenzidine	119-93-7	7.30E-04	c	6.8E-04	c	7.30E-04	NA	NA	3.00E+00	T	3.00E+00
2-Acetylaminofluorene	53-96-3	NA	NA	NA	NA	NA	NA	NA	2.50E+03	T	2.50E+03
bis(2-Ethylhexyl)phthalate	1117-81-7	4.80E-01	c	4.47E-01	c	4.80E-01	NA	NA	1.00E+04	T	1.00E+04
3,3'-Dichlorobenzidine	91-94-1	1.50E-02	c	1.4E-02	c	1.50E-02	NA	NA	6.21E+03	T	6.21E+03
Benz(a)anthracene	56-55-3	2.20E-02	c	8.6E-03	c	2.20E-02	NA	NA	6.00E+02	T	6.00E+02
Chrysene	2118-01-9	2.17E+00	c	8.58E-01	c	2.17E+00	NA	NA	2.00E+02	T	2.00E+02
Oi-n-octylphthalate	1117-84-0	7.30E+01	nc	7.30E+01	nc	7.30E+01	NA	NA	1.50E+05	T	1.50E+05
7,12-Dimethylbenz(a)anthracene	57-97-6	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzot(b)fluoranthene	205-99-2	2.20E-02	c	8.6E-03	c	2.20E-02	NA	NA	NA	NA	NA
Benzot(k)fluoranthene	207-08-9	2.20E-01	c	8.6E-02	c	2.20E-01	NA	NA	NA	NA	NA
Benz(a)pyrene	50-32-8	2.20E-03	c	2.0E-03	c	2.20E-03	NA	NA	7.50E+03	T	7.50E+03
3-Methylcholanthrene	56-49-5	NA	NA	NA	NA	NA	NA	NA	1.50E+03	T	1.50E+03
Indeno(1,2,3-cd)pyrene	193-39-5	2.17E-02	c	8.58E-03	c	2.17E-02	NA	NA	NA	NA	NA
Dibenz(a,h)anthracene	53-70-3	2.17E-03	c	8.58E-04	c	2.17E-03	NA	NA	3.00E+04	T	3.00E+04
Benzo(g,h,i)perylene	191-24-2	NA	NA	NA	NA	NA	NA	NA	3.00E+04	T	3.00E+04

Footnotes:

PRG: Preliminary Remediation Goals

c: Cancer

nc: non-cancer

RBC: Risk-Based Concentration

HBSL: Health-based Screening Level

(E) ERPG: Emergency Response Planning Guidelines

(T) TEEL: Temporary Emergency Exposure Limits

ATV: Acute Toxicity Value

NA: Not available

APPENDIX D

RISK EVALUATION DATA

Table D-1: Comparison of Air Concentrations With Health-Based Values: Metals, Particulates and Miscellaneous Compounds

Compound	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	$C_{\text{chronic}}/\text{HBSL}$	Red Parachute Signal Flare			$C_{\text{acute}}/\text{ATV}$	$>1?$
				$>1?$	$C_{\text{acute}} (\mu\text{g}/\text{m}^3)$	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)		
TSP	1.78E-02	5.00E+01	3.55E-04	no	NA	NA	na	na
PM ₁₀	1.68E-02	5.00E+01	3.36E-04	no	NA	NA	na	na
HCl (a)	NA	2.08E+01	4.97E-05	no	NA	7.14E+03	na	na
Cl ₂ (a)	1.04E-05	2.09E-01	7.25E-07	no	1.82E-02	2.89E+03	6.28E-06	no
Dioxin TEQ (b)	3.25E-14	4.48E-08	9.03E-06	no	5.31E-10	3.50E+00	1.52E-10	no
Carbon Monoxide (CO)	1.42E-03	1.57E+02	9.17E-06	no	2.49E+00	2.30E+05	1.08E-05	no
Nitrogen Oxide (NOx)	4.17E-04	1.00E+02	4.17E-06	no	2.92E+00	2.70E+05	1.08E-05	no
HCl (a)	1.42E-05	2.08E+01	6.80E-07	no	9.93E-02	7.14E+03	1.39E-05	no
Carbon Dioxide (CO ₂)	1.94E-02	NV	na	1.36E+02	5.40E+07	2.52E-06	no	no
Sulfur Dioxide (SO ₂)	9.84E-06	8.00E+01	1.23E-07	no	1.72E-02	7.89E+02	2.18E-05	no
Aluminum	2.32E-05	3.65E+00	6.36E-06	no	1.63E-01	3.00E+04	5.42E-06	no
Antimony	6.07E-08	1.46E+00	4.16E-08	no	4.26E-04	1.50E+03	2.84E-07	no
Arsenic	NA	4.47E-04	na	NA	3.00E+01	na	na	na
Barium	4.91E-05	5.21E-01	9.43E-05	no	3.44E-01	1.50E+03	2.30E-04	no
Beryllium	NA	8.00E-04	na	NA	5.00E+00	na	na	na
Cadmium	3.85E-08	1.07E-03	3.61E-05	no	6.30E-04	3.00E+01	2.10E-05	no
Chromium	2.00E-07	1.53E-04	1.31E-03	no	3.27E-03	1.50E+03	2.18E-06	no
Cobalt	6.10E-08	2.20E-02	2.77E-10	no	4.27E-04	6.00E+01	7.12E-06	no
Copper	9.40E-07	1.46E-02	6.44E-09	no	6.59E-03	3.00E+03	2.20E-06	no
Lead	3.44E-07	1.50E-00	2.29E-07	no	2.41E-03	1.50E+02	1.61E-05	no
Magnesium	2.12E-03	NV	na	1.49E+01	3.00E+04	4.95E-04	no	no
Manganese	3.23E-06	5.11E-02	6.32E-05	no	2.26E-02	3.00E+03	7.55E-06	no
Nickel	9.35E-08	7.30E-01	1.28E-09	no	6.55E-04	3.00E+03	2.18E-07	no
Phosphorus	1.60E-06	NV	na	1.12E-02	3.00E+02	3.73E-05	no	no
Selenium	NA	1.83E+01	na	NA	6.00E+02	na	na	na
Silver	NA	1.83E+01	na	NA	3.00E+02	na	na	na
Thallium	NA	2.56E-01	na	NA	3.00E+02	na	na	na
Zinc	8.93E-07	1.10E+03	8.15E-10	no	6.26E-03	3.00E+04	2.09E-07	no
Mercury	1.27E-14	3.13E-01	4.07E-14	no	8.92E-05	1.00E+02	8.92E-07	no

Footnote:

(a) HCl/Cl₂ levels were too low to be reliably measured.

(b) Presence questionable - reported at similar levels in samples and blanks.
NA = Not applicable because compound was not detected.

na = Not available because health-based screening value is not available or not applicable if compound was not detected.
NV = No value
 C_{chronic} = Chronic time-averaged concentration ; HBSL = Chronic health-based screening level

$>1?$ = Is the ratio greater than one?
 C_{acute} = Acute concentration; ATV = Acute toxicity value

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	Red Parachute Signal Flare						> 1?
	C _{chronic} (µg/m ³)	Health-Based Screening Level (µg/m ³)	C _{chronic} /HBSL	> 1?	C _{acute} (µg/m ³)	Acute Toxicity Value (µg/m ³)	
Total Nonmethane Hydrocarbons (TNMHC)							
TNMC	4.55E-05	NV		na	NA	NA	na
Volatile Organic Compounds (VOCs)							
Ethane	1.68E-06	NV		na	NA	NA	na
Ethylene	1.76E-05	NV		na	1.23E-01	4.60E+05	2.68E-07
Acetylene	1.22E-05	NV		na	NA	NA	na
Propane	3.78E-07	NV		na	2.65E-03	3.78E+06	7.01E-10
Propene	3.86E-06	NV		na	NA	NA	na
i-Butane	7.56E-08	NV		na	5.30E-04	5.71E+06	9.28E-11
i-Butene	2.12E-07	NV		na	NA	NA	na
1-Butene	7.56E-07	NV		na	NA	NA	na
1,3-Butadiene	4.15E-07	3.74E-03	1.11E-04	no	1.70E-03	2.20E+04	7.71E-08
n-Butane	1.06E-07	NV		na	7.42E-04	5.71E+06	1.30E-10
trans-2-Butene	7.56E-07	NV		na	NA	NA	na
2,2-Dimethylpropane	NA	NV		na	NA	NA	na
cis-2-Butene	1.36E-07	NV		na	NA	NA	na
3-Methyl-1-butene	NA	NV		na	NA	NA	na
i-Pentane	NA	NV		na	NA	1.80E+06	na
1-Pentene	1.51E-07	NV		na	NA	NA	na
2-Methyl-1-butene	3.02E-08	NV		na	NA	NA	na
n-Pentane	1.81E-07	NV		na	1.27E-03	1.80E+06	7.07E-10
Isoprene	1.51E-07	NV		na	NA	NA	na
trans-2-Pentene	3.02E-08	NV		na	NA	NA	na
cis-2-Pentene	3.02E-08	NV		na	NA	NA	na
2-Methyl-2-butene	6.05E-08	NV		na	NA	NA	na
2,2-Dimethylbutane	6.05E-08	NV		na	4.24E-04	1.80E+06	2.36E-10
Cyclopentene	4.54E-08	NV		na	NA	NA	na
4-Methyl-1-pentene	NA	NV		na	NA	NA	na
Cyclopentane	1.51E-08	NV		na	NA	NA	na

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	Red Parachute Signal Flare					
	C _{chronic} (µg/m ³)	Health-Based Screening Level (µg/m ³)	C _{chronic} /HBSL	> 1?	C _{acute} (µg/m ³)	Acute Toxicity Value (µg/m ³)
2,3-Dimethylbutane	1.51E-08	NV		na	NA	NA
cis-4-Methyl-2-pentene	NA	NV		na	NA	na
2-Methylpentane	3.02E-08	NV		na	2.12E-04	1.80E+06
3-Methylpentane	NA	NV		na	NA	1.18E-10
2-Methyl-1-pentene	NA	NV		na	NA	no
1-Hexene	1.21E-07	NV		na	NA	na
n-Hexane	3.02E-08	2.10E+02	1.44E-10	no	8.48E-04	1.03E+05
trans-2-Hexene	NA	NV		na	2.12E-04	8.23E-09
2-Methyl-2-pentene	NA	NV		na	NA	no
cis-2-Hexene	NA	NV		na	NA	na
Methylcyclopentane	4.54E-08	NV		na	NA	NA
2,4-Dimethylpentane	NA	NV		na	NA	na
Benzene	7.52E-07	2.50E-01	3.01E-06	no	3.07E-03	1.56E+05
Cyclohexane	1.21E-07	NV		na	8.48E-04	3.10E+06
2-Methylhexane	NA	NV		na	NA	2.74E-10
2,3-Dimethylpentane	NA	NV		na	NA	no
3-Methylhexane	NA	NV		na	NA	na
2,2,4-Trimethylpentane	NA	NV		na	NA	na
n-Heptane	7.56E-08	NV		na	5.30E-04	1.97E-08
2,4,4-Trimethyl-1-pentene	NA	NV		na	NA	no
Methylcyclohexane	NA	3.10E+03		na	NA	na
2,4,4-Trimethyl-2-pentene	NA	NV		na	NA	na
2,5-Dimethylhexane	1.51E-08	NV		na	NA	na
2,4-Dimethylhexane	NA	NV		na	NA	na
2,3,4-Trimethylpentane	1.64E-23	NV		na	NA	NA
Toluene	3.63E-07	4.02E+02	9.04E-10	no	6.36E-04	1.88E+05
2,3-Dimethylhexane	NA	NV		na	NA	3.39E-09
2-Methylheptane	3.02E-08	NV		na	NA	na
3-Ethylhexane	3.02E-08	NV		na	NA	na

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	C _{chronic} (µg/m ³)	Health-Based Screening Level (µg/m ³)	C _{chronic/HBSL}	Red Parachute Signal Flare				
				>1?	C _{acute} (µg/m ³)	Acute Toxicity Value (µg/m ³)	C _{acute/ATV}	>1?
2,2-Dimethylheptane	NA	NV		na	NA	NA		na
2,2,4-Trimethylhexane	NA	NV		na	NA	NA		na
n-Octane	3.02E-08	NV		na	NA	NA		na
Ethylcyclohexane	NA	NV		na	NA	NA		na
Ethylbenzene	NA	1.10E+03		na	NA	5.43E+05		na
m-Xylene & p-Xylene	NA	NV		na	NA	6.51E+05		na
Styrene	1.36E-07	1.10E+03	1.24E-10	no	2.38E-04	2.13E+05	1.12E-09	no
o-Xylene	NA	7.30E+03		na	NA	6.51E+05		na
n-Nonane	2.27E-07	4.02E+02	5.65E-10	no	1.59E-03	1.05E+06	1.52E-09	no
i-Propylbenzene	NA	4.00E+02		na	NA			na
n-Propylbenzene	6.05E-08	3.65E+01	1.66E-09	no	NA			na
p-Ethyltoluene	1.51E-07	NV		na	1.06E-03	1.25E+05	8.48E-09	no
m-Ethyltoluene	9.07E-08	NV		na	NA			na
1,3,5-Trimethylbenzene	7.56E-08	6.20E+00	1.22E-08	no	5.30E-04	3.68E+05	1.44E-09	no
o-Ethyltoluene	1.21E-07	NV		na	8.48E-04	7.50E+02	1.13E-06	no
1,2,4-Trimethylbenzene & sec-Butylbenzene	1.81E-07	6.21E+00	2.92E-08	no	1.27E-03	1.80E+05	7.07E-09	no
n-Decane	NA	NV		na	NA	4.37E+03		na
alpha-Pinene	NA	NV		na	NA	4.00E+04		na
beta-Pinene	NA	NV		na	NA	NA		na
delta 3-Carene	NA	NV		na	NA			na
d-Limonene	NA	NV		na	NA	1.95E+06		na
MTBE	NA	3.10E+03		na	NA	4.32E+05		na
Dichlorodifluoromethane	NA	2.10E+02		na	NA	1.48E+07		na
Methylchloride	NA	NV		na	NA			na
Dichlorotetrafluoroethane	NA	NV		na	NA			na
Chloroethene	NA	2.20E-02		na	NA	1.28E+04		na
1,3-Butadiene	4.22E-07	3.74E-03	1.13E-04	no	1.72E-03	2.20E+04	7.84E-08	no
Methylbromide	NA	5.20E+00		na	NA	5.82E+04		na

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	Red Parachute Signal Flare						>1?
	C _{chronic} (µg/m ³)	Health-Based Screening Level (µg/m ³)	C _{chronic} /HBSL	>1?	C _{acute} (µg/m ³)	Acute Toxicity Value (µg/m ³)	
Ethylchloride	1.81E-08	2.30E+00	7.88E-09	no	2.96E-04	7.92E+06	3.74E-11
Trichloromonofluoromethane	NA	7.30E+02		na	NA	2.81E+06	no
Vinyldenechloride	NA	NV		na	NA	7.92E+04	na
Methylenechloride	NA	4.10E+00		na	NA	6.96E+05	na
Allylchloride	NA	1.00E+00		na	NA	9.39E+03	na
1,1,2-Trichloro-1,2,2-trifluoroethane	NA	3.13E+04		na	NA	9.58E+06	na
1,1-Dichloroethane	NA	5.21E+02		na	NA	1.21E+06	na
1,2-Dichloroethene	NA	3.29E+01		na	NA	2.38E+06	na
Chloroform	NA	8.40E-02		na	NA	9.76E+03	na
1,2-Dichloroethane	NA	7.39E-02		na	NA	8.08E+03	na
Methylchloroform	NA	1.00E+03		na	NA	1.91E+06	na
Benzene	7.65E-07	2.50E-01	3.06E-06	no	1.25E-02	1.60E+05	7.84E-08
Carbontetrachloride	NA	1.04E+03		na	NA	1.28E+05	na
1,2-Dichloropropane	NA	9.89E-02		na	NA	5.08E+05	na
Trichloroethylene	NA	1.12E+00		na	NA	5.37E+05	na
cis 1,3-Dichloro-1-propene	NA	NV		na	NA	1.14E+04	na
trans 1,3-Dichloro-1-propene	NA	NV		na	NA	NA	na
1,1,2-Trichloroethane	NA	1.20E-01		na	NA	1.64E+05	na
Toluene	3.69E-07	4.02E+02	9.19E-10	no	6.47E-04	1.88E+05	3.45E-09
1,2-Dibromoethane	NA	8.73E-03		na	NA	1.54E+05	na
Perchloroethylene	NA	3.31E+00		na	NA	6.89E+05	na
Chlorobenzene	NA	6.20E+01		na	NA	1.38E+05	na
Ethylbenzene	NA	1.10E+03		na	NA	4.34E+03	na
m&p-Xylene	NA	7.30E+02		na	NA	6.51E+05	na
Styrene	1.38E-07	1.06E+03	1.31E-10	no	2.43E-04	2.13E+05	1.14E-09
1,1,2,2-Tetrachloroethane	NA	3.31E-02		na	NA	2.06E+04	na
o-Xylene	NA	7.30E+02		na	NA	6.51E+05	na
p-Ethyltoluene	1.54E-07	NV		na	1.08E-03	1.25E+05	8.62E-09
1,3,5-Trimethylbenzene	9.23E-08	6.21E+00	1.49E-08	no	6.47E-04	3.68E+05	1.76E-09

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	Red Parachute Signal Flare						C _{acute} /ATV	>1?
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	C _{chronic} /HBSL	>1?	C _{acute} ($\mu\text{g}/\text{m}^3$)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)		
1,2,4-Trimethylbenzene	1.85E-07	6.21E+00	2.97E-08	no	1.29E-03	1.80E+05	7.19E-09	no
Benzylchloride	NA	4.00E-02		na	NA	5.20E+03		na
m-Dichlorobenzene	NA	3.30E+00		na	NA	3.61E+04		na
p-Dichlorobenzene	NA	2.80E-01		na	NA	6.61E+05		na
o-Dichlorobenzene	NA	2.09E+02		na	NA	3.01E+05		na
1,2,4-Trichlorobenzene	NA	NV		na	NA	3.71E+04		na
Hexachlorobutadiene	NA	8.73E-02		na	NA	3.21E+04		na
trans-1,2-Dichloroethene	NA	7.30E+01		na	NA	4.95E+04		na
o-Chlorotoluene	NA	7.30E+01		na	NA	3.88E+05		na
p-Chlorotoluene	NA	NV		na	NA	3.88E+05		na
1,3,5-Trichlorobenzene	NA	NV		na	NA	NA		na
1,2,3-Trichlorobenzene	NA	NV		na	NA	5.00E+04		na
Methylnitrite	3.15E-08	NV		na	NA	NA		na
Acetonitrile	1.51E-07	6.20E+01	2.43E-09	no	1.05E-03	1.01E+05	1.05E-08	no
Acrylonitrile	4.94E-08	2.80E-02	1.77E-06	no	2.02E-04	2.20E+04	9.18E-09	no
Nitromethane	7.49E-08	NV		na	5.25E-04	1.50E+05	3.50E-09	no
Benzonitrile	6.50E-08	NV		na	4.55E-04	1.50E+04	3.04E-08	no
Nitrobenzene	NA	2.09E+00		na	NA	1.51E+04		na
Carbonyl Sulfide	6.22E-08	NV		na	4.36E-04	9.84E+03	4.43E-08	no
Sulfur Dioxide	NA	NV		na	NA	7.80E+02		na
Carbon Disulfide	1.86E-06	7.30E+02	2.55E-09	no	1.31E-02	3.73E+04	3.50E-07	no
Thiophene	7.81E-08	NV		na	NA	NA		na
Dimethyldisulfide	NA	NV		na	NA	4.00E+01		na
2-Methylthiophene	NA	NV		na	NA	NA		na
3-Methylthiophene	NA	NV		na	NA	NA		na
Dimethyltrisulfide	NA	NV		na	NA	NA		na
Isothiocyanatomethane	NA	NV		na	NA	NA		na
2-Chlorothiophene	NA	NV		na	NA	NA		na
3-Chlorothiophene	NA	NV		na	NA	NA		na

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	C _{chronic} (µg/m ³)	Health-Based Screening Level (µg/m ³)	C _{chronic} /HBSL	Red Parachute Signal Flare			
				> 1?	C _{acute} (µg/m ³)	Acute Toxicity Value (µg/m ³)	C _{acute} / ATV > 1?
2-Thiophenecarboxaldehyde	NA	NV		na	NA	NA	na
Naphthalene	9.90E-08	3.13E+00	3.17E-08	no	6.94E-04	7.86E+04	8.83E-09 no
Acetaldehyde	8.24E-08	8.70E-01	9.47E-08	no	3.37E-04	1.80E+04	1.87E-08 no
Acrolein	8.01E-08	2.10E-02	3.82E-06	no	1.40E-04	2.30E+02	6.10E-07 no
Acetone	NA	3.40E+02		na	NA	2.37E+06	na
Propanal	1.82E-07	NV		na	1.27E-03	7.50E+04	1.70E-08 no
Furan	7.99E-08	3.70E+00	2.16E-08	no	5.60E-04	1.67E+02	3.36E-06 no
2-Propanol	NA	NV		na	NA	9.84E+05	na
2-Methylpropanal	NA	NV		na	NA	NA	na
Methacrolein	NA	NV		na	NA	NA	na
2,3-Butanedione	NA	NV		na	NA	NA	na
Methyl-Vinyl Ketone	NA	NV		na	NA	8.61E+01	na
MTBE	1.22E-08	3.10E+03	3.93E-12	no	8.54E-05	4.32E+05	1.98E-10 no
Butanal	8.97E-08	NV		na	6.29E-04	7.38E+04	8.52E-09 no
2-Butanone	1.42E-07	1.00E+03	1.42E-10	no	9.93E-04	8.85E+05	1.12E-09 no
Tetrahydrofuran	NA	9.89E-01		na	NA	7.38E+05	na
2-Methyl-1-propanol	NA	1.10E+03		na	NA	4.55E+05	na
trans-2-Butenal	NA	3.54E-03		na	NA	NA	na
Acetic Acid	1.52E-07	NV		na	1.07E-03	3.68E+04	2.90E-08 no
2-Pentanone	1.76E-07	NV		na	1.23E-03	8.80E+05	1.40E-09 no
Pentanal	1.42E-08	NV		na	NA	NA	na
4-Methyl-2-pentanone	NA	8.30E+01		na	NA	3.07E+05	na
trans-2-Pentenal	NA	NV		na	NA	NA	na
Cyclopentanone	NA	NV		na	NA	NA	na
2-Hexanone	6.60E-08	5.11E+00	1.29E-08	no	4.62E-04	4.09E+04	1.13E-08 no
Hexanal	NA	NV		na	NA	NA	na
3-Furaldehyde	NA	NV		na	NA	NA	na
Butyl Acetate	NA	NV		na	NA	NA	na
2-Furaldehyde	NA	5.20E+01		na	NA	7.86E+03	na

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	Red Parachute Signal Flare							
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	C _{chronic} /HBSL	> 1?	C _{acute} ($\mu\text{g}/\text{m}^3$)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	C _{acute} / ATv	> 1?
trans-2-Hexenal	NA	NV		na	NA	NA	na	na
1-Hexanol	NA	NV		na	NA	8.36E+03	na	na
3-Heptanone	1.52E-07	NV		na	NA	NA	na	na
2-Heptanone	NA	NV		na	NA	1.70E+03	na	na
Heptanal	NA	NV		na	NA	NA	na	na
trans-2-Heptenal	NA	NV		na	NA	NA	na	na
5-Methyl-2-furaldehyde	NA	NV		na	NA	NA	na	na
6-Methyl-2-heptanone	NA	NV		na	NA	NA	na	na
Benzaldehyde	7.10E-08	3.70E+02	1.92E-10	no	4.97E-04	1.50E+04	3.32E-08	no
1-Heptanol	NA	NV		na	NA	NA	na	na
6-Methyl-5-hepten-2-one	NA	NV		na	NA	NA	na	na
2-Octanone	NA	NV		na	NA	NA	na	na
Octanal	NA	NV		na	NA	NA	na	na
Benzofuran	NA	NV		na	NA	NA	na	na
trans-2-Octenal	NA	NV		na	NA	NA	na	na
Acetophenone	2.65E-08	2.10E-02	1.26E-06	no	1.86E-04	3.00E+04	6.20E-09	no
2-Nonanone	NA	NV		na	NA	NA	na	na
Nonanal	5.41E-07	NV		na	NA	NA	na	na
trans-2-Nonenal	NA	NV		na	NA	NA	na	na
2-Decanone	NA	NV		na	NA	NA	na	na
Decanal	8.05E-07	NV		na	NA	NA	na	na

Footnotes:

(a) Items in bold represent duplicate values for those compounds that are common for Method TO-14 and TO-12.

NA = Not applicable

na = Not available because health-based screening value is not available or not applicable because compound was not detected.

NV = No value

C_{chronic} = Chronic time-averaged concentration

HBSL = Chronic health-based screening level

>? = Is the ratio greater than one?

Table D-2: Comparison of Air Concentrations With Health-Based Values: Volatile Organic Compounds

Compound (a)	Red Parachute Signal Flare							
	C_{chronic} ($\mu\text{g}/\text{m}^3$)	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	$C_{\text{chronic}}/\text{HBSL}$	$> 1?$	C_{acute} ($\mu\text{g}/\text{m}^3$)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	$C_{\text{acute}}/\text{ATV}$	$> 1?$

C_{acute} = Acute concentration

ATV = Acute toxicity value

Table D-3: Comparison of Air Concentrations With Health-Based Values: Semi-Volatile Organic Compounds

Compound	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	C _{chronic} / HBSL	Red Parachute Signal Flare			
				> 1?	C _{acute} ($\mu\text{g}/\text{m}^3$)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	C _{acute} / ATV
Particulate/Vapor-phase SVOCs							
N-Nitrosodimethylamine	NA	1.40E-04		na	NA	2.50E+03	na
Pyridine	NA	3.65E+00		na	NA	4.85E+04	na
2-Picoline	NA	NV		na	NA	NA	na
Methyl methanesulfonate	NA	NV		na	NA	NA	na
N-Nitrosomethylalkylamine	NA	3.06E-04		na	NA	NA	na
N-Nitrosodiethylamine	NA	4.47E-05		na	NA	NA	na
Ethyl methanesulfonate	NA	NV		na	NA	NA	na
Phenol	NA	2.19E+03		na	NA	3.85E+05	na
Aniline	NA	1.06E+00		na	NA	2.29E+04	na
bis(2-Chloroethyl)ether	NA	5.80E-03		na	NA	5.85E+04	na
Pentachloroethane	NA	NV		na	NA	NA	na
2-Chlorophenol	NA	1.80E+01		na	NA	5.25E+03	na
1,3-Dichlorobenzene	NA	NV		na	NA	NA	na
1,4-Dichlorobenzene	NA	2.80E-01		na	NA	6.61E+05	na
Benzyl alcohol	NA	1.10E+03		na	NA	5.53E+04	na
2-Methylphenol	NA	NV		na	NA	6.63E+04	na
1,2-Dichlorobenzene	NA	2.09E+02		na	NA	3.01E+05	na
bis(2-Chloroisopropyl)ether	NA	1.92E-01		na	NA	6.99E+04	na
o-Toluidine	NA	2.80E-02		na	NA	2.63E+04	na
4-Methylphenol/3-Methylphenol	NA	NV		na	NA	6.63E+04	na
N-Nitroso-di-n-propylamine	NA	9.61E-04		na	NA	5.32E+03	na
Acetophenone	3.14E-07	2.10E-02	1.50E-05	no	2.20E-03	1.47E+05	1.50E-08
N-Nitrosomorpholine	NA	NV		na	NA	3.00E+04	na
N-Nitrosopyrrolidine	NA	3.15E-03		na	NA	NA	na
Hexachloroethane	NA	4.80E-01		na	NA	2.90E+04	na
Nitrobenzene	NA	2.09E+00		na	NA	1.51E+04	na
N-Nitrosopiperidine	NA	NV		na	NA	NA	na
Isophorone	NA	7.08E+00		na	NA	2.83E+04	na
2,4-Dimethylphenol	NA	7.30E+01		na	NA	NA	na
2-Nitrophenol	NA	NV		na	NA	NA	na

Table D-3: Comparison of Air Concentrations With Health-Based Values: Semi-Volatile Organic Compounds

Compound	Red Parachute Signal Flare						> 1?
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	C _{chronic} / HBSL	> 1?	C _{acute} ($\mu\text{g}/\text{m}^3$)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	
bis(2-Chloroethoxy)methane	NA	NV			NA	NA	na
Benzoic acid	3.56E-07	1.50E+04	2.37E-11	no	2.49E-03	1.25E+04	1.99E-07
2,4-Dichlorophenol	NA	1.10E+01		na	NA	3.00E+04	na
1,2,4-Trichlorobenzene	NA	NV		na	NA	3.71E+04	na
Naphthalene	NA	3.13E+00		na	NA	7.86E+04	na
p-Chloroaniline	NA	1.46E+01		na	NA	5.21E+03	na
2,6-Dichlorophenol	NA	NV		na	NA	3.00E+04	na
Hexachloropropene	NA	NV		na	NA	NA	na
Hexachlorobutadiene	NA	8.73E-02		na	NA	3.21E+04	na
Dimethylphenethylamine	NA	NV		na	NA	NA	na
N-Nitroso-di-n-butylamine	NA	1.20E-03		na	NA	NA	na
4-Chloro-3-methylphenol	NA	NV		na	NA	NA	na
Safrole	NA	NV		na	NA	NA	na
2-Methylnaphthalene	NA	NV		na	NA	2.00E+04	na
1,2,4,5-Tetrachlorobenzene	NA	1.10E+00		na	NA	3.00E+04	na
Hexachlorocyclopentadiene	NA	7.30E-02		na	NA	2.23E+02	na
2,4,6-Trichlorophenol	NA	6.20E-01		na	NA	3.00E+04	na
2,4,5-Trichlorophenol	NA	3.70E+02		na	NA	3.00E+04	na
Isosafrole	NA	NV		na	NA	NA	na
2-Chloronaphthalene	NA	2.90E+02		na	NA	6.00E+02	na
2-Nitroaniline	NA	2.10E-01		na	NA	NA	na
1,4-Naphthoquinone	NA	NV		na	NA	2.50E+02	na
Dimethylphthalate	NA	3.65E+04		na	NA	1.50E+04	na
1,3-Dinitrobenzene	NA	3.70E-01		na	NA	3.00E+03	na
2,6-Dinitrotoluene	NA	3.70E+00		na	NA	6.00E+02	na
Acenaphthylene	NA	NV		na	NA	2.00E+02	na
3-Nitroaniline	NA	NV		na	NA	NA	na
4-Nitrophenol	NA	2.90E+01		na	NA	3.00E+04	na
2,4-Dinitrophenol	NA	7.30E+00		na	NA	7.50E+03	na.
Acenaphthene	NA	2.20E+02		na	NA	1.25E+03	na
2,4-Dinitrotoluene	NA	7.30E+00		na	NA	6.00E+02	na

Table D-3: Comparison of Air Concentrations With Health-Based Values: Semi-Volatile Organic Compounds

Compound	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	C _{chronic/ HBSL}	> 1?	Red Parachute Signal Flare		
					C _{acute} ($\mu\text{g}/\text{m}^3$)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	C _{acute/} ATV
Dibenzofuran	NA	1.46E+01	na	NA	1.50E+00	na	na
Pentachlorobenzene	NA	2.92E+00	na	NA	3.00E+04	na	na
1-Naphthylamine	NA	NV	na	NA	3.50E+04	na	na
2-Naphthylamine	NA	NV	na	NA	7.50E+03	na	na
2,3,4,6-Tetrachlorophenol	NA	1.10E+02	na	NA	NA	na	na
Diethylphthalate	4.55E-08	2.92E+03	1.56E-11	no	3.19E-04	1.50E+04	2.13E-08
4-Chlorophenylphenyl ether	NA	NV	na	NA	NA	na	na
Fluorene	NA	1.46E+02	na	NA	7.50E+04	na	na
5-Nitro-o-toluidine	NA	2.00E-01	na	NA	NA	na	na
4-Nitroaniline	NA	NV	na	NA	9.00E+03	na	na
4,6-Dinitro-2-methylphenol	NA	3.65E-01	na	NA	5.00E+02	na	na
Diphenylamine/N-NitrosodPAA	NA	NV	na	NA	2.50E+03	na	na
sym-Trinitrobenzene	NA	1.10E+02	na	NA	3.00E+04	na	na
Diallate	NA	1.10E-01	na	NA	NA	na	na
Phenacetin	NA	NV	na	NA	3.00E+04	na	na
4-Bromophenylphenyl ether	NA	NV	na	NA	NA	na	na
Hexachlorobenzene	NA	4.18E-03	na	NA	7.50E+01	na	na
4-Aminobiphenyl	NA	NV	na	NA	1.50E+03	na	na
Pronamide	NA	2.74E+02	na	NA	NA	na	na
Pentachlorophenol	NA	5.60E-02	na	NA	1.50E+03	na	na
Pentachloronitrobenzene	NA	2.59E-02	na	NA	1.50E+03	na	na
Phenanthrene	NA	NV	na	NA	2.00E+03	na	na
Anthracene	NA	1.10E+03	na	NA	6.00E+03	na	na
Carbazole	NA	3.36E-01	na	NA	NA	na	na
Di-n-butylphthalate	NA	3.65E+02	na	NA	1.50E+04	na	na
4-Nitroquinoline-1-oxide	NA	NV	na	NA	NA	na	na
Methapyrilene	NA	NV	na	NA	NA	na	na
Fluoranthene	NA	1.50E+02	na	NA	3.00E+01	na	na
Benzidine	NA	2.90E-05	na	NA	5.00E+02	na	na
Pyrene	NA	NV	na	NA	1.50E+04	na	na
p-Dimethylaminobenzene	NA	NV	na	NA	7.50E+04	na	na

Table D-3: Comparison of Air Concentrations With Health-Based Values: Semi-Volatile Organic Compounds

Compound	Red Parachute Signal Flare							
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Health-Based Screening Level ($\mu\text{g}/\text{m}^3$)	C _{chronic} / HBSL	> 1?	C _{acute} ($\mu\text{g}/\text{m}^3$)	Acute Toxicity Value ($\mu\text{g}/\text{m}^3$)	C _{acute} / ATV	> 1?
Chlorobenzilate	NA	2.49E-02		na	NA	2.50E+02		na
Kepone	NA	3.74E-04		na	NA	1.00E+02		na
Butylbenzylphthalate	1.12E-07	7.30E+02	1.53E-10	no	7.83E-04	5.00E+05	1.57E-09	no
3,3'-Dimethylbenzidine	NA	7.30E-04		na	NA	3.00E+00		na
2-Acetylaminofluorene	NA	NV		na	NA	2.50E+03		na
bis(2-Ethylhexyl)phthalate	3.10E-08	4.80E-01	6.45E-08	no	5.07E-04	1.00E+04	5.07E-08	no
3,3'-Dichlorobenzidine	NA	1.50E-02		na	NA	6.21E+03		na
Benz(a)anthracene	NA	2.20E-02		na	NA	6.00E+02		na
Chrysene	NA	2.17E+00		na	NA	2.00E+02		na
Di-n-octylphthalate	6.68E-08	7.30E+01	9.15E-10	no	4.68E-04	1.50E+05	3.12E-09	no
7,12-Dimethylbenz(a)anthracene	NA	NV		na	NA	NA		na
Benzo(b)fluoranthene	NA	2.20E-02		na	NA	NA		na
Benzo(k)fluoranthene	NA	2.20E-01		na	NA	NA		na
Benz(a)pyrene	NA	2.20E-03		na	NA	7.50E+03		na
3-Methylcholanthrene	NA	NV		na	NA	1.50E+03		na
Indeno(1,2,3-cd)pyrene	NA	2.17E-02		na	NA	NA		na
Dibenz(a,h)anthracene	NA	2.17E-03		na	NA	3.00E+04		na
Benzog(h,i)perylene	NA	NV		na	NA	3.00E+04		na

Footnotes:

NA = Not applicable

na = Not available because health-based screening value is not available or not applicable because compound was not detected.

NV = No value

C_{chronic} = Chronic time-averaged concentration

HBSL = Chronic health-based screening level

>1? = Is the ratio greater than one?

C_{acute} = Acute concentration

ATV = Acute toxicity value

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)			
TNMHC	NA	NA	NA	NA
Ethane	NA	NA	NA	NA
Ethylene	NA	NA	NA	NA
Acetylene	NA	NA	NA	NA
Propane	3.78E-07	NA	NA	NA
Propene	3.86E-06	NA	NA	NA
i-Butane	7.56E-08	NA	NA	NA
i-Butene	2.12E-07	NA	NA	NA
1-Butene	7.56E-07	NA	NA	NA
1,3-Butadiene	NA	NA	NA	NA
n-Butane	1.06E-07	NA	NA	NA
trans-2-Butene	7.56E-07	NA	NA	NA
2,2-Dimethylpropane	NA	NA	NA	NA
cis-2-Butene	1.36E-07	NA	NA	NA
3-Methyl-1-butene	NA	NA	NA	NA
i-Pentane	NA	NA	NA	NA
1-Pentene	1.51E-07	NA	NA	NA
2-Methyl-1-butene	3.02E-08	NA	NA	NA
n-Pentane	1.81E-07	NA	NA	NA
Isoprene	NA	NA	NA	NA
trans-2-Pentene	3.02E-08	NA	NA	NA
cis-2-Pentene	3.02E-08	NA	NA	NA
2-Methyl-2-butene	6.05E-08	NA	NA	NA
2,2-Dimethylbutane	6.05E-08	NA	NA	NA
Cyclopentene	4.54E-08	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)			
	<i>Aliphatic:C<=8</i>	<i>Aliphatic:C>8</i>	<i>Aromatic:C<=8</i>	<i>Aromatic:C>8</i>
4-Methyl-1-pentene	NA	NA	NA	NA
Cyclopentane	1.51E-08	NA	NA	NA
2,3-Dimethylbutane	1.51E-08	NA	NA	NA
cis-4-Methyl-2-pentene	NA	NA	NA	NA
2-Methylpentane	3.02E-08	NA	NA	NA
3-Methylpentane	NA	NA	NA	NA
2-Methyl-1-pentene	NA	NA	NA	NA
1-Hexene	1.21E-07	NA	NA	NA
n-Hexane	3.02E-08	NA	NA	NA
trans-2-Hexene	NA	NA	NA	NA
2-Methyl-2-pentene	NA	NA	NA	NA
cis-2-Hexene	NA	NA	NA	NA
Methylcyclopentane	4.54E-08	NA	NA	NA
2,4-Dimethylpentane	NA	NA	NA	NA
Benzene	NA	NA	1.75E-06	NA
Cyclohexane	1.21E-07	NA	NA	NA
2-Methylhexane	NA	NA	NA	NA
2,3-Dimethylpentane	NA	NA	NA	NA
3-Methylhexane	NA	NA	NA	NA
2,2,4-Trimethylpentane	NA	NA	NA	NA
n-Heptane	7.56E-08	NA	NA	NA
2,4,4-Trimethyl-1-pentene	NA	NA	NA	NA
Methylcyclohexane	NA	NA	NA	NA
2,4,4-Trimethyl-2-pentene	NA	NA	NA	NA
2,5-Dimethylhexane	1.51E-08	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)			
	<i>Aliphatic:C<=8</i>	<i>Aliphatic:C>8</i>	<i>Aromatic:C<=8</i>	<i>Aromatic:C>8</i>
2,4-Dimethylhexane	NA	NA	NA	NA
2,3,4-Trimethylpentane	1.64E-23	NA	NA	NA
Toluene	NA	NA	3.63E-07	NA
2,3-Dimethylhexane	NA	NA	NA	NA
2-Methylheptane	3.02E-08	NA	NA	NA
3-Ethylhexane	3.02E-08	NA	NA	NA
2,2-Dimethylheptane	NA	NA	NA	NA
2,2,4-Trimethylhexane	NA	NA	NA	NA
n-Octane	3.02E-08	NA	NA	NA
Ethylcyclohexane	NA	NA	NA	NA
Ethylbenzene	NA	NA	NA	NA
m-Xylene & p-Xylene	NA	NA	NA	NA
Styrene	NA	NA	NA	1.36E-07
o-Xylene	NA	NA	NA	NA
n-Nonane	NA	2.27E-07	NA	NA
i-Propylbenzene	NA	NA	NA	NA
n-Propylbenzene	NA	NA	NA	6.05E-08
p-Ethyltoluene	NA	NA	NA	1.51E-07
m-Ethyltoluene	NA	NA	NA	9.07E-08
1,3,5-Trimethylbenzene	NA	NA	NA	7.56E-08
o-Ethyltoluene	NA	NA	NA	1.21E-07
1,2,4-Trimethylbenzene & sec-Butylbenzene	NA	NA	NA	1.81E-07
n-Decane	NA	NA	NA	NA
alpha-Pinene	NA	NA	NA	NA
beta-Pinene	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)			
Alliphatic:C<=8		Alliphatic:C>8		Aromatic:C<=8
delta 3-Carene	NA	NA	NA	NA
d-Limonene	NA	NA	NA	NA
MTBE	NA	NA	NA	NA
Dichlorodifluoromethane	NA	NA	NA	NA
Methylchloride	NA	NA	NA	NA
Dichlorotetrafluoroethane	NA	NA	NA	NA
Chloroethene	NA	NA	NA	NA
1,3-Butadiene	NA	NA	NA	NA
Methylbromide	NA	NA	NA	NA
Ethylchloride	NA	NA	NA	NA
Trichloromonofluoromethane	NA	NA	NA	NA
Vinyldenechloride	NA	NA	NA	NA
Methylenechloride	NA	NA	NA	NA
Allylchloride	NA	NA	NA	NA
1,1,2-Trichloro-1,2,2-trifluoroethane	NA	NA	NA	NA
1,1-Dichloroethane	NA	NA	NA	NA
1,2-Dichloroethene	NA	NA	NA	NA
Chloroform	NA	NA	NA	NA
1,2-Dichloroethane	NA	NA	NA	NA
Methylchloroform	NA	NA	NA	NA
Benzene	NA	NA	1.78E-06	NA
Carbontetrachloride	NA	NA	NA	NA
1,2-Dichloropropane	NA	NA	NA	NA
Trichloroethylene	NA	NA	NA	NA
cis 1,3-Dichloro-1-propene	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare				Aromatic:C>8
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Aromatic:C<=8	C _{chronic} ($\mu\text{g}/\text{m}^3$)	
trans 1,3-Dichloro-1-propene	NA	NA	Aliphatic:C<=8	NA	NA
1,1,2-Trichloroethane	NA	NA	Aliphatic:C>8	NA	NA
Toluene	NA	NA	Aliphatic:C<=8	NA	NA
1,2-Dibromoethane	NA	NA	Aliphatic:C>8	3.69E-07	NA
Perchloroethylene	NA	NA	Aliphatic:C<=8	NA	NA
Chlorobenzene	NA	NA	Aliphatic:C>8	NA	NA
Ethylbenzene	NA	NA	Aliphatic:C<=8	NA	NA
m&p-Xylene	NA	NA	Aliphatic:C>8	NA	NA
Styrene	NA	NA	Aliphatic:C<=8	NA	1.38E-07
1,1,2,2-Tetrachloroethane	NA	NA	Aliphatic:C>8	NA	NA
o-Xylene	NA	NA	Aliphatic:C<=8	NA	NA
p-Ethyltoluene	NA	NA	Aliphatic:C>8	NA	1.54E-07
1,3,5-Trimethylbenzene	NA	NA	Aliphatic:C<=8	NA	9.23E-08
1,2,4-Trimethylbenzene	NA	NA	Aliphatic:C>8	NA	1.85E-07
Benzylchloride	NA	NA	Aliphatic:C<=8	NA	NA
m-Dichlorobenzene	NA	NA	Aliphatic:C>8	NA	NA
p-Dichlorobenzene	NA	NA	Aliphatic:C<=8	NA	NA
o-Dichlorobenzene	NA	NA	Aliphatic:C>8	NA	NA
1,2,4-Trichlorobenzene	NA	NA	Aliphatic:C<=8	NA	NA
Hexachlorobutadiene	NA	NA	Aliphatic:C>8	NA	NA
trans-1,2-Dichloroethene	NA	NA	Aliphatic:C<=8	NA	NA
o-Chlorotoluene	NA	NA	Aliphatic:C>8	NA	NA
p-Chlorotoluene	NA	NA	Aliphatic:C<=8	NA	NA
1,3,5-Trichlorobenzene	NA	NA	Aliphatic:C>8	NA	NA
1,2,3-Trichlorobenzene	NA	NA	Aliphatic:C<=8	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)			
Methylnitrite	NA	NA	NA	NA
Acetonitrile	NA	NA	NA	NA
Acrylonitrile	NA	NA	NA	NA
Nitromethane	NA	NA	NA	NA
Benzonitrile	NA	NA	NA	NA
Nitrobenzene	NA	NA	NA	NA
Carbonyl Sulfide	NA	NA	NA	NA
Sulfur Dioxide	NA	NA	NA	NA
Carbon Disulfide	NA	NA	NA	NA
Thiophene	NA	NA	NA	NA
Dimethyldisulfide	NA	NA	NA	NA
2-Methylthiophene	NA	NA	NA	NA
3-Methylthiophene	NA	NA	NA	NA
Dimethyltrisulfide	NA	NA	NA	NA
Isothiocyanatomethane	NA	NA	NA	NA
2-Chlorothiophene	NA	NA	NA	NA
3-Chlorothiophene	NA	NA	NA	NA
2-Thiocphenecarboxaldehyde	NA	NA	NA	NA
Naphthalene	NA	NA	NA	9.90E-08
Acetaldehyde	NA	NA	NA	NA
Acrolein	NA	NA	NA	NA
Acetone	NA	NA	NA	NA
Propanal	NA	NA	NA	NA
Furan	NA	NA	NA	NA
2-Propanol	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Aromatic:C>8	Aromatic:C<=8
2-Methylpropanal	NA	NA	NA	NA
Methacrolein	NA	NA	NA	NA
2,3-Butanedione	NA	NA	NA	NA
Methyl-Vinyl Ketone	NA	NA	NA	NA
MTBE	NA	NA	NA	NA
Butanal	NA	NA	NA	NA
2-Butanone	NA	NA	NA	NA
Tetrahydrofuran	NA	NA	NA	NA
2-Methyl-1-propanol	NA	NA	NA	NA
trans-2-Butenal	NA	NA	NA	NA
Acetic Acid	NA	NA	NA	NA
2-Pentanone	NA	NA	NA	NA
Pentanal	NA	NA	NA	NA
4-Methyl-2-pentanone	NA	NA	NA	NA
trans-2-Pentenal	NA	NA	NA	NA
Cyclopentanone	NA	NA	NA	NA
2-Hexanone	NA	NA	NA	NA
Hexanal	NA	NA	NA	NA
3-Furaldehyde	NA	NA	NA	NA
Butyl Acetate	NA	NA	NA	NA
2-Furaldehyde	NA	NA	NA	NA
trans-2-Hexenal	NA	NA	NA	NA
1-Hexanol	NA	NA	NA	NA
3-Heptanone	NA	NA	NA	NA
2-Heptanone	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Aromatic:C<=8	Aromatic:C>8
Heptanal	NA	NA	NA	NA
trans-2-Heptenal	NA	NA	NA	NA
5-Methyl-2-furaldehyde	NA	NA	NA	NA
6-Methyl-2-heptanone	NA	NA	NA	NA
Benzaldehyde	NA	NA	NA	NA
1-Heptanol	NA	NA	NA	NA
6-Methyl-5-hepten-2-one	NA	NA	NA	NA
2-Octanone	NA	NA	NA	NA
Octanal	NA	NA	NA	NA
Benzofuran	NA	NA	NA	NA
trans-2-Octenal	NA	NA	NA	NA
Acetophenone	NA	NA	NA	NA
2-Nonanone	NA	NA	NA	NA
Nonanal	NA	NA	NA	NA
trans-2-Nonenal	NA	NA	NA	NA
2-Decanone	NA	NA	NA	NA
Decanal	NA	NA	NA	NA
N-Nitrosodimethylamine	NA	NA	NA	NA
Pyridine	NA	NA	NA	NA
2-Picoline	NA	NA	NA	NA
Methyl methanesulfonate	NA	NA	NA	NA
N-Nitrosomethylethyamine	NA	NA	NA	NA
N-Nitrosodiethylamine	NA	NA	NA	NA
Ethyl methanesulfonate	NA	NA	NA	NA
Phenol	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$
Aniline	NA	NA	NA	NA
bis(2-Chloroethyl)ether	NA	NA	NA	NA
Pentachloroethane	NA	NA	NA	NA
2-Chlorophenol	NA	NA	NA	NA
1,3-Dichlorobenzene	NA	NA	NA	NA
1,4-Dichlorobenzene	NA	NA	NA	NA
Benzyl alcohol	NA	NA	NA	NA
2-Methylphenol	NA	NA	NA	NA
1,2-Dichlorobenzene	NA	NA	NA	NA
bis(2-Chloroisopropyl)ether	NA	NA	NA	NA
o-Toluidine	NA	NA	NA	NA
4-Methylphenol/3-Methylphenol	NA	NA	NA	NA
N-Nitroso-di-n-propylamine	NA	NA	NA	NA
Acetophenone	NA	NA	NA	NA
N-Nitrosomorpholine	NA	NA	NA	NA
N-Nitrosopyrrolidine	NA	NA	NA	NA
Hexachloroethane	NA	NA	NA	NA
Nitrobenzene	NA	NA	NA	NA
N-Nitrosopiperidine	NA	NA	NA	NA
Isophorone	NA	NA	NA	NA
2,4-Dimethylphenol	NA	NA	NA	NA
2-Nitrophenol	NA	NA	NA	NA
bis(2-Chloroethoxy)methane	NA	NA	NA	NA
Benzoic acid	NA	NA	NA	NA
2,4-Dichlorophenol	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)	C _{chronic} ($\mu\text{g}/\text{m}^3$)	Aromatic:C<=8	Aromatic:C>8
1,2,4-Trichlorobenzene	NA	NA	NA	NA
Naphthalene	NA	NA	NA	NA
p-Chloroaniline	NA	NA	NA	NA
2,6-Dichlorophenol	NA	NA	NA	NA
Hexachloropropene	NA	NA	NA	NA
Hexachlorobutadiene	NA	NA	NA	NA
Dimethylphenethylamine	NA	NA	NA	NA
N-Nitroso-di-n-butylamine	NA	NA	NA	NA
4-Chloro-3-methylphenol	NA	NA	NA	NA
Safrole	NA	NA	NA	NA
2-Methylnaphthalene	NA	NA	NA	NA
1,2,4,5-Tetrachlorobenzene	NA	NA	NA	NA
Hexachlorocyclopentadiene	NA	NA	NA	NA
2,4,6-Trichlorophenol	NA	NA	NA	NA
2,4,5-Trichlorophenol	NA	NA	NA	NA
Isosafrole	NA	NA	NA	NA
2-Chloronaphthalene	NA	NA	NA	NA
2-Nitroaniline	NA	NA	NA	NA
1,4-Naphthoquinone	NA	NA	NA	NA
Dimethylphthalate	NA	NA	NA	NA
1,3-Dinitrobenzene	NA	NA	NA	NA
2,6-Dinitrotoluene	NA	NA	NA	NA
Acenaphthylene	NA	NA	NA	NA
3-Nitroaniline	NA	NA	NA	NA
4-Nitrophenol	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$
2,4-Dinitrophenol	NA	NA	NA	NA
Acenaphthene	NA	NA	NA	NA
2,4-Dinitrotoluene	NA	NA	NA	NA
Dibenzo furan	NA	NA	NA	NA
Pentachlorobenzene	NA	NA	NA	NA
1-Naphthylamine	NA	NA	NA	NA
2-Naphthylamine	NA	NA	NA	NA
2,3,4,6-Tetrachlorophenol	NA	NA	NA	NA
Diethyl phthalate	NA	NA	NA	NA
4-Chlorophenyl phenyl ether	NA	NA	NA	NA
Fluorene	NA	NA	NA	NA
5-Nitro-o-tolidine	NA	NA	NA	NA
4-Nitroaniline	NA	NA	NA	NA
4,6-Dinitro-2-methylphenol	NA	NA	NA	NA
Diphenylamine/N-NitrosodPA	NA	NA	NA	NA
sym-Trinitrobenzene	NA	NA	NA	NA
Diallate	NA	NA	NA	NA
Phenacetin	NA	NA	NA	NA
4-Bromophenyl phenyl ether	NA	NA	NA	NA
Hexachlorobenzene	NA	NA	NA	NA
4-Aminobiphenyl	NA	NA	NA	NA
Pronamide	NA	NA	NA	NA
Pentachlorophenol	NA	NA	NA	NA
Pentachloronitrobenzene	NA	NA	NA	NA
Phenanthrene	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	C _{chronic} ($\mu\text{g}/\text{m}^3$)			
Anthracene	Aliphatic:C<=8 NA	Aliphatic:C>8 NA	Aromatic:C<=8 NA	Aromatic:C>8 NA
Carbazole	NA	NA	NA	NA
Di-n-butylphthalate	NA	NA	NA	NA
4-Nitroquinoline-1-oxide	NA	NA	NA	NA
Methapryllene	NA	NA	NA	NA
Fluoranthene	NA	NA	NA	NA
Benzidine	NA	NA	NA	NA
Pyrene	NA	NA	NA	NA
p-Dimethylaminoazobenzene	NA	NA	NA	NA
Chlorobenzilate	NA	NA	NA	NA
Kepone	NA	NA	NA	NA
Butylbenzylphthalate	NA	NA	NA	NA
3,3'-Dimethylbenzidine	NA	NA	NA	NA
2-Acetylaminofluorene	NA	NA	NA	NA
bis(2-Ethylhexyl)phthalate	NA	NA	NA	NA
3,3'-Dichlorobenzidine	NA	NA	NA	NA
Benz(a)anthracene	NA	NA	NA	NA
Chrysene	NA	NA	NA	NA
Di-n-octylphthalate	NA	NA	NA	NA
7,12-Dimethylbenz(a)anthracene	NA	NA	NA	NA
Benzo(b)fluoranthene	NA	NA	NA	NA
Benzo(k)fluoranthene	NA	NA	NA	NA
Benz(a)pyrene	NA	NA	NA	NA
3-Methylcholanthrene	NA	NA	NA	NA
Indeno(1,2,3-cd)pyrene	NA	NA	NA	NA

Table D-4: Comparison of Air Concentrations With Health-Based Values: Total Petroleum Hydrocarbons

Compound (a)	Red Parachute Signal Flare			
	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$	$C_{\text{chronic}} (\mu\text{g}/\text{m}^3)$
	$A\text{liphatic:}C\leq 8$	$A\text{liphatic:}C>8$	$A\text{romatic:}C\leq 8$	$A\text{romatic:}C>8$
Dibenz(a,h)anthracene	NA	NA	NA	NA
Benzo(g,h)perylene	NA	NA	NA	NA
Total ($\mu\text{g}/\text{m}^3$)	7.43E-06	2.27E-07	2.15E-06	7.87E-07
Derived Health-Based Screening Level	1.92E+04	1.04E+03	4.17E+02	2.09E+02
$C_{\text{chronic}}/\text{HBSL}$	3.87E-10	2.18E-10	5.16E-09	3.77E-09
>1?	no	no	no	no

Footnotes:

(a) Items in bold represent duplicate values. highest concentration was used to estimate total petroleum hydrocarbon concentration
 >1? = Is the ratio greater than one?

NA = Not Applicable because compound was not detected

C_{chronic} = Chronic averaged air Concentration

HBSL = Health-Based Screening Level

APPENDIX E

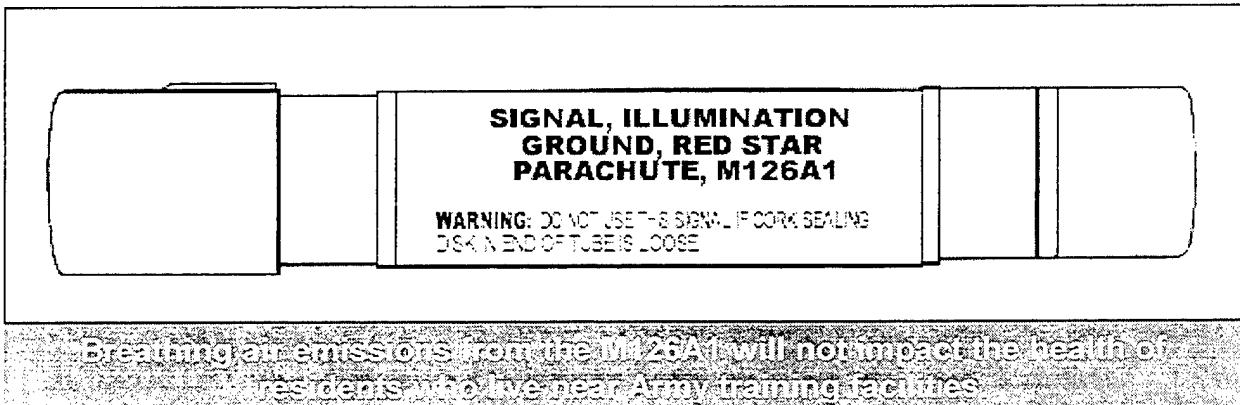
FACT SHEET SUBMITTED TO AEC

United States Army Environmental Center

Pyrotechnics Fact Sheet

M126A1 Red Star Parachute Signal Flare

Department of Defense Identification Code: L311



WHAT ARE PYROTECHNICS?

The terms pyrotechnics and fireworks are often used interchangeably. Pyrotechnics are devices that give off smoke, light, and/or a loud noise when activated. Military pyrotechnics are used for signaling, obscuring, and illuminating during training and combat.

WHAT IS THE M126A1?

The M126A1 is a star parachute signal flare, which is a type of pyrotechnic device used for signaling and illuminating. The M126A1 produces a single, red, parachute-suspended illuminating star. It is 10.16 inches long, 1.67 inches wide, and weighs 1.20 pounds.

HOW IS THE M126A1 USED?

A rocket containing the signal is launched from a hand-held device. After igniting, the rocket reaches a height of

about 200 feet and produces a single, red star illumination resembling a firework. The signal extends to a height of 700 to 750 feet and can be seen from a distance of 30 to 35 miles at night. Use of this device is important in training our troops to use and identify different signals, which is an important method of communication in the field.

WHERE IS THE M126A1 USED?

The M126A1 is used during many Army training events. These events are held at nearly every Army training installation. At most locations, the training areas are at least 1000 meters (over half a mile) away from populated areas. In general, one item is used during a day of training, which typically occurs five times per year.

WHAT IS IN THE M126A1?

The M126A1 consists of a parachute-suspended illumination assembly and a rocket motor propulsion assembly. These

are contained in a hand-held aluminum launching tube. The illumination component consists primarily of strontium nitrate, and magnesium powder.

WILL BREATHING AIR EMISSIONS FROM THE M126A1 AFFECT MY HEALTH?

To answer this question, the U.S. Army Environmental Center tested the air emissions from the M126A1. The U.S. Army Center for Health Promotion and Preventive Medicine then determined if there would be a potential for health effects from inhalation to residents living near training areas. Results showed that residents breathing air as close as 100 meters (328 feet) from the activation site are safe from these emissions.

HOW WAS THE STUDY DONE?

To gather data for the study, airborne emissions were collected by activating the M126A1 in a test chamber. The air in the chamber was tested to identify the types and amount of substances released. More than 300 substances were looked for during this part of the study.

This information was then used in an air model (a computer program that allows estimation of air concentrations) to determine the amount of each substance, to which someone living near a training area might be exposed. Downwind concentrations were estimated based on a typical use

scenario for the M126A1. Since the study does not look at a specific training area, the assumptions used in the model will in most cases, predict higher downwind air concentrations than those expected at an actual training site.

These estimated air concentrations were then compared to safe screening levels established by the U.S. Environmental Protection Agency and other agencies. If the air concentrations are below these screening levels, they are considered safe for everyone, including sensitive people such as the sick, elderly, and children.

WHAT ARE THE LIMITATIONS OF THIS STUDY?

Many steps were taken to ensure that the results of this study are protective of everyone who lives close to training areas. However, limitations do exist with this study. For example, the study does not consider exposure to other types of munitions that could also be used during the same training event. Due to these limitations, conservative model conditions were used to ensure the protection of public health from inhalation of the M126A1 air emissions.

WHERE CAN I GET MORE INFORMATION?

For more information on the M126A1 and other military munitions call the Army Environmental Center Hotline at 1-800-USA-3845, visit our website at www.aec.army.mil, or email us at t2hotline@aec.apgea.army.mil.